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**Supplemental Ground Water  
Investigation Report  
I-70, 44<sup>th</sup> Street to Brighton  
Boulevard, City and County of  
Denver, Colorado**

WALSH Project Number: 3026-010  
CDOT Project No. IR-IM (CX) 070-4 (145)  
September 15, 1999

**ADMINISTRATIVE RECORD**



Environmental Scientists and Engineers, Inc.

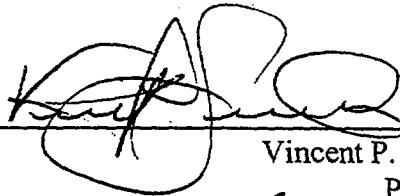
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# **SUPPLEMENTAL GROUND WATER INVESTIGATION REPORT, I-70, 44<sup>TH</sup> STREET TO BRIGHTON BOULEVARD, CITY AND COUNTY OF DENVER, COLORADO**

September 15, 1999

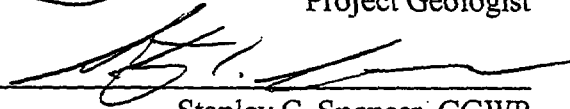
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WALSH Project Number: 3026-010

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## EXECUTIVE SUMMARY

The findings of this investigation indicate that ground water containing PCE is entering the CDOT project area near I-70 and Brighton Boulevard from the southeast, and flowing under the project area toward the northwest. The plume of PCE contamination appears to extend from at least as far south as 44<sup>th</sup> Street and Brighton Boulevard, to at least as far northeast as the 4600 block of Brighton Boulevard, north of I-70 and Brighton Boulevard.

The concentrations of PCE were highest between TH-24 (530 µg/L and 480 µg/L in 1998 and 1999, respectively), and TH-26 (160 µg/L in 1999). PCE concentrations at TH-25 (23 µg/L in 1999) and at a number of locations near Brighton Boulevard and 44<sup>th</sup> Street suggest that concentrations may decrease to the northeast and southwest of the areas of highest observed PCE concentrations.

The findings of this investigation indicate ground water containing PCE at concentrations exceeding Colorado Ground Water Standards is entering the CDOT project area from upgradient sources, and that the PCE plume has extended under numerous properties. The full extent and origin of the PCE plume have not been delineated.

WALSH recommends that CDOT transmit the findings of this investigation to the Colorado Department of Public Health and Environment (CDPHE). CDPHE or perhaps the U.S. EPA may wish to conduct further investigations to determine the origin and extent of the PCE contaminated ground water in this area. The PCE contaminated ground water appears to originate from sources outside of CDOT's project area.

WALSH recommends that CDOT consider that any ground water which may be encountered during construction in the vicinity of I-70 and Brighton Boulevard may contain PCE in excess of Colorado Ground Water Standards. Any construction activities that involve handling of ground water will require prior testing of ground water for PCE, and proper handling and possibly treatment of PCE contaminated ground water.

The U.S. EPA is conducting an investigation of arsenic and lead in residential soils in this area of Denver. In this study, arsenic and lead concentrations were measured in surface soil at three non-residential locations. The highest observed arsenic concentration (38 mg/Kg at TH-26) was below target levels of the U.S. EPA investigation of residential soils in this area. The highest observed lead concentration (850 mg/Kg at TH-26) exceeded the target concentration of the U.S. EPA investigation of residential soils in this area. WALSH recommends that CDOT transmit the arsenic and lead data for these surface soils to CDPHE and the U.S. EPA.

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# **SUPPLEMENTAL GROUND WATER INVESTIGATION REPORT I-70, 44<sup>TH</sup> STREET TO BRIGHTON BOULEVARD, CITY AND COUNTY OF DENVER**

## **INTRODUCTION**

Walsh Environmental Scientists and Engineers, Inc. (WALSH) has conducted a supplemental ground water investigation in the study area as requested by the Colorado Department of Transportation (CDOT). Tetrachloroethene (PCE) was discovered in ground water during Phase II Site Investigation activities for CDOT Project No. IR-IM(CX)070-4(145). CDOT's construction project involves widening of I-70 from Washington Street to High Street in Denver, Colorado. The PCE contamination was discovered during previous investigations on CDOT's construction project (WALSH, 1991a, 1991b, 1992, 1996 and 1998).

The purpose of this 1999 investigation was to determine whether previously identified PCE contamination in ground water originated within the area of CDOT's construction project, or from sources upgradient of CDOT's project area. Certain additional laboratory analyses were performed on some soil and ground water samples collected during this investigation. Lead and arsenic were measured in surface soils at three locations at the request of the Colorado Department of Health and Environment. Volatile organic compounds (VOCs), including benzene, toluene, ethylbenzene and xylenes (BTEX), total extractable petroleum hydrocarbons (EPA), and semivolatile organic compounds (SVOCs) were measured in soils from three test areas to provide additional information on the distribution of various potential contaminants in the study area.

CDOT is in the process of converting to metric system. Therefore this document is in metric measurement format with American equivalents following in parenthesis, where appropriate. Where a manufacturer's unit is given (e.g., 1,000 gallon tank), or a unit is common to professional literature, or is quoted from other sources, the original unit is retained.

The units milligrams per kilogram (mg/Kg) and milligrams per liter are used in place of parts per million (ppm) for solids and liquids, respectively, throughout this document. Similarly, micrograms per kilogram ( $\mu\text{g/Kg}$ ) and micrograms per liter ( $\mu\text{g/L}$ ) are used in place of parts per billion (ppb). However, measurements of headspace gases are reported in ppm.

The figures referenced in the body of this document are presented in Appendix A. Tables are presented in Appendix B, soil boring/monitoring well construction details in Appendix C, field data transmittal sheets in Appendix D, and laboratory data in Appendix E.

## PROPOSED CONSTRUCTION AND PROPERTY ACQUISITION

The CDOT is replacing the elevated portion of I-70 from Washington Street to Brighton Boulevard with a wider, elevated highway (see Figure 2). Existing ramps at Humboldt Street will be removed and replaced with interchanges constructed at Washington Street and Brighton Boulevard. The existing elevated structure east of Humboldt Street will be replaced, and fill material will be used to support I-70 east of Humboldt Street.

Design plans indicate that property acquisition is required for the widening of I-70, the widening of Washington Street, and the relocations of 46th Avenue and Brighton Boulevard. The Union Pacific Railroad (UPRR) tracks (south of I-70 at Brighton Boulevard) will be moved southward to accommodate proposed ramp structures. Construction of retaining walls, ditches, storm and sanitary sewers, and other utility relocations will require excavation at various locations in the project area.

## REGIONAL SETTING

### 1 Location

The area investigated is located within the City and County of Denver, Colorado and is shown in Figure 1. The initial phase of the construction will involve the widening of the eastbound lanes of I-70 from Washington Street to Humboldt Street. Properties affected by the Phase I and Phase II construction activities were evaluated in WALSH, 1997 and 1998. Properties included in this investigation are located between 44<sup>th</sup> Street and Brighton Boulevard, south of I-70. The properties to be sampled have been or will be acquired by CDOT. Construction activities will extend ramps away from I-70 and into existing neighborhoods and commercial areas approximately bounded by 44th Street and Brighton Boulevard on the south and 47th Avenue on the north. The UPRR tracks will be relocated south of their present position.

### 2 Physiography and Geology

The project area is located in the Denver Basin, east of the Front Range Uplift of the Southern Rocky Mountains. Topographically the area is generally flat with elevations decreasing towards the South Platte River (Figure 1). Surface drainage is towards the South Platte River. The project area is covered by a thin veneer of unconsolidated sediments, including the Roadway Alluvium (Quaternary) and the Post Piney Creek Alluvium (Holocene). These sediments are generally poorly sorted sands, gravel, and some clays. Flat lying bedrock of the Cretaceous/Tertiary Denver Formation, consisting of weathered shale, siltstone, and fine sandstone, is unconformably overlain by the younger unconsolidated sediments. Depth to the bedrock surface near I-70 and east of Humboldt Street is approximately 45 feet (Guirre, 1995).



The top of the Denver Formation marks the base of the unconfined aquifer. Depth to ground water is approximately 27 to 34 feet below ground surface in the vicinity of Humboldt Street and Brighton Boulevard (WALSH, 1991b, 1992 and Aguirre, 1995). Local, unconfined ground water flow east of the South Platte River is to the northwest (WALSH, 1991b). Ground water flow velocities were estimated to vary between 20 and 200 feet per year within the study area (WALSH, 1996). More details of the local geology and ground water conditions are found in the various WALSH reports (1991b, 1992, 1996, 1997, and 1998) and the geotechnical report by Aguirre (1995).

## **REVIEW OF EARLIER ENVIRONMENTAL INVESTIGATIONS**

Phase II and III of the modifications to I-70 between Humboldt Street and Brighton Boulevard traverse a commercial, industrial and residential area where many environmental concerns have been identified. Several properties of concern were identified between 44<sup>th</sup> Street and Brighton Boulevard in 1991 and 1992 (WALSH, 1991a, 1991b, and 1992), including nine properties with known or suspected USTs. This study was expanded in July 1991 to include the I-70 corridor from Washington Street to Brighton Boulevard and was revised in late 1996. These reports identified four main categories of environmental concern: petroleum contaminated soils and ground water from leaking USTs and ASTs; possible soil and ground water contamination from tannery operations; and soil contaminated with smelter wastes resulting in elevated heavy metal content. Chlorinated solvents (PCE) were discovered in several monitoring wells and sampling performed during March of 1999 has confirmed the presence of PCE at concentrations above safe drinking water levels. However no source has been positively identified for these contaminants. Abandoned USTs of unknown use were removed from three locations during late 1998 and early 1999 (WALSH, 1999). Laboratory analyses performed during the tank removals did not detect chlorinated solvents in either the tank contents or surrounding soils. Analyses detected only diesel range hydrocarbons which indicated that the tanks had probably stored fuel oil.

This investigation was intended to determine if PCE in ground water originated within CDOT's project area, or migrated onto CDOT ROW from unknown up gradient sources. This work also provided confirmation of the presence of PCE within the project area, and additional information on the extent of the PCE contamination.

## **SITE ASSESSMENT ACTIVITIES**

WALSH drilled three additional test holes/monitor wells (TH-26, TH-27, and TH-28) on June 7 and June 18, 1999 in the southeast (up gradient) area of the project to identify possible offsite migration onto the property from the southeast. The location of ground water test holes TH-26, TH-27, TH-28 are shown on Figure 2.

WALSH sampled the three newly installed test holes, and re-sampled four previously existing test holes TH-19, TH-25, TH-24, and TH-7 (W) on March 23, 1999, to help determine the down gradient extent of contamination and possible impacts to offsite receptors.

Ground water was analyzed for volatile organic compounds (VOCs) and total extractable petroleum hydrocarbons (TEPH). Soil samples from drilling activities were also analyzed for Cs and TEPH in addition to total lead and arsenic and semi-volatile organic compounds (SVOCs) from drill cuttings of surface soils.

Static water levels were surveyed in the wells to refine potentiometric surface maps and help determine ground water flow directions and pathways.

As part of an ongoing Corrective Action Plan and Final Site Characterization for a former underground storage tank (UST) at Parcel 49, WALSH sampled ground water at temporary monitor locations PZ-1, PZ-2 and PZ-3 (See Figure 2) on June 25, 1999. Ground water samples at PZ-1, PZ-2 and PZ-3 were also analyzed for PCE contamination. The sampling analysis of ground water at PZ-1, PZ-2 and PZ-3 are presented in Corrective Action Plan Final Site Characterization Report Parcel 49, Tank 2, I-70 and Brighton Boulevard, Denver, Colorado (WALSH, 1999).

## **SAMPLING METHODOLOGIES**

### **Soil Sampling**

The locations of the test holes are shown on Figure 2. Test holes were drilled to a depth of approximately 9.1 meters (30 feet) below ground surface (bgs). Test holes were drilled using each hollow stem continuous flight augers. Test holes were monitored during drilling for flammable combustible gases and VOCs using a calibrated combustible gas indicator (CGI) and a calibrated photoionization detector (PID). Sample headspace measurements were taken using PID to field screen for VOCs.

Soil samples from all test holes were collected from drill cuttings from the ground surface interval to approximately 0 to 0.9 meters (0-3 feet) bgs and thereafter, at 5-foot intervals using stainless-steel split spoon sampler. The soil sample from the 0 to 0.9 meter (0-3 feet), and additional soil samples collected from the most suspect interval and soil/ground water surface zone were sent to the laboratory. All samples were collected in glass jars with non-leaked lids. Samples were kept in coolers for delivery to the laboratory. Chain-of-custody records were completed for each sample. Complete logs were prepared for all test holes.

### **Ground Water Sampling**

Ground water was sampled through factory-slotted (screened) PVC pipe inserted in the test holes. Screened PVC sections were installed from total depth to at least 2 feet above the static water table. Solid 2-inch PVC sections with a locking cap was used to extend the screened interval to the surface. Silica sand was poured around the PVC to approximately 2 feet above

screen, and bentonite chips were added from the sand level to approximately 1.5 feet below the surface to seal the annular space. Concrete was used to secure a flush mount. Test holes (TH-26, TH-27, TH-28) were developed by purging a minimum of 10 casing volumes of water. Remaining test holes to be sampled were purged of three casing volumes. Prior to sample collection, the well headspace was measured for VOCs using a PID and the standing water was field tested during bailing for pH, conductivity and temperature. The stability of these measurements and the total volume of water purged assured that the ground water sample is representative of the formation. Ground water sampling forms were completed for each test hole.

Ground water samples were retrieved with a disposable polypropylene bailer and collected for analysis as follows: two 40-milliliter (ml) VOA vials for VOCs; and one 1-liter (L) glass bottles for TEPH. All samples were placed immediately in an ice-filled cooler and delivered to the laboratory with complete chain-of-custody records.

### 3 Surveying

The potentiometric surface of this area is known from previous studies (WALSH 1991b, 1998). The ground level and casing elevations of newly installed test holes was recorded by a global positioning system (GPS) by Drexell Barrell, Inc. of Colorado Springs, using the dual frequency method, as specified in the contract with CDOT.

## LABORATORY ANALYSES

Soil and ground water samples collected for semi-volatile organic compounds (SVOCs, EPA Method 8270), volatile organic compounds (VOCs, EPA Method 8260), and total extractable petroleum hydrocarbons (TEPH, diesel range hydrocarbons, EPA Modified Method 8100), were analyzed at the WALSH Laboratory in Boulder, Colorado. Soil samples collected for total lead and arsenic (EPA Methods SW 7421 and 7060) were analyzed at Analytical Laboratories, Inc., in Broomfield, Colorado. Tables 3, 4, and 5 summarize the analytical results for soil, and Table 6 summarizes the analytical results for ground water.

## FIELD OBSERVATIONS

### 1 Test Hole Drilling

Field observations during drilling activities indicated similar lithology for the three test holes. From 0 to 0.9 meters (0-3 feet), black, fill material (coal dust) consisting of silty clay was encountered. Lithology from 1.5 meters (5 feet) to 6.1 meters (20 feet) consisted of coarse to medium sand with layers of silty sand and intermittent layers of clayey silt and sand present. The water/ground water interface zone at roughly 7.6 meters (25 feet) and phreatic zone at 9.1 meters (30 feet), consisted predominantly of coarse to fine sand and medium to fine gravel.

Staining and odors were not observed throughout the profile of each test hole, however, PID readings were registered. PID readings from TH-26 ranged from 0 to 7 ppm. PID readings from TH-27 ranged from 0.1 ppm to 23.6 ppm. PID readings from TH-28 ranged from 0 ppm to 1.6 ppm.

The soil ground water interface zone was encountered during drilling activities at roughly the 7.6 to 8.2 meters (25-27 feet) bgs interval. At 9.1 to 9.8 meters (30-32 feet) bgs, the soil was completely saturated.

## 8.2 Ground Water Sampling

Ground water samples were collected from pre-existing test holes TH-7 (W), TH-19, TH-24, and TH-25 on March 23, 1999. Test holes TH-26, TH-27, and TH-28 were developed and sampled on June 23, 1999.

Ground water in TH-26, TH-27, and TH-28 was silty (brown) in appearance at the onset of development, and remained silty at the time of sampling. Sheen, odors, discoloration were not observed in ground water.

## 10 ANALYTICAL RESULTS

### 10.1 Ground Water

#### 10.1.1 VOCs, TVPH and TEPH

Concentrations of PCE in ground water exceeded Colorado Ground Water Standards at all sample locations. Concentrations of PCE in ground water for down gradient test holes sampled on March 23, 1999 indicated the following concentrations: TH-7 (W), 61 µg/L; TH-19, 160 µg/L; TH-24, 480 µg/L; TH-25, 23 µg/L.

Concentrations of PCE in ground water for up gradient test holes sampled on June 23, 1999 indicated the following concentrations: TH-26, 300 µg/L; TH-27, 340 µg/L; TH-28, 160 µg/L (See Figure 3). All concentrations exceed the maximum contaminant level of 5 µg/L for PCE in ground water.

As reported separately (WASLH, 1999) PCE was found in ground water at PZ-1 (44 µg/L), PZ-2 (33 µg/L) and PZ-3 (7.8 µg/L). No ground water was collected at PZ-4.

TVPH and TEPH concentrations were non-detectable for all test holes, with the exception of a TEPH concentration in ground water at TH-7 (W) of 1,300 µg/L, and a TVPH concentration in ground water at TH-24 of 600 µg/L. Currently, there are no adopted state ground water standards for TVPH and TEPH.

## 0.2 Soil

### 0.2.1 VOCs and TEPH

Analytical results for soil samples collected for analyses of VOCs (BTEX included) indicated non-detectable concentrations for all three newly installed test holes TH-26, TH-27, and TH-28. A TEPH concentration of 11 mg/Kg was detected in TH-26 at 0 to 0.3 meters (0-1 feet) bgs, and at 6 mg/Kg from 3 to 3.7 meters (10-12 feet) in soil at TH-28. These concentrations of TEPH, however, are in compliance with Department of Labor and Employment, Oil Inspection Section (DOLE-OIS) Tier 1 Risk Based Screening Levels (RBSLs) of 500 mg/Kg (DOLE-OIS, 1999). Table 3 in Appendix B summarizes analytical results for VOCs and TEPH.

### 0.2.2 SVOCs

Analytical results for shallow soil samples collected for analyses of SVOCs indicated non-detectable concentrations for TH-26 and TH-27. Detectable levels of phenanthrene (0.34 mg/Kg), fluoranthene (0.56 mg/Kg), pyrene (0.92 mg/Kg), chrysene (0.37 mg/Kg), and benzo(b)fluoranthene (0.43 mg/Kg) were indicated in soil from 0.3 to 0.9 meters (1-3 feet) bgs at TH-28. These concentrations, however, are below residential and industrial surficial soil Tier 1 RBSLs. Table 4 in Appendix B summarizes analytical results for SVOCs.

### 0.2.3 Total Arsenic and Lead

Analytical results for shallow soil samples collected for analyses of total arsenic and lead indicated detectable levels in TH-26, TH-27, and TH-28. The soil sample from 0 to 0.3 meters (0-1 feet) bgs at TH-26 contained a total arsenic concentration of 38 mg/Kg, and total lead of 850 mg/Kg. The soil sample from 0 to 0.3 meters (0-1 feet) bgs at TH-27 contained a total arsenic concentration of 4.7 mg/Kg, and total lead of 68 mg/Kg. The soil sample from 0.3 to 0.9 meters (1-3 feet) bgs contained a total arsenic concentration of 15 mg/Kg, and total lead of 140 mg/Kg. These detectable levels are consistent with background metal concentrations typically associated with soils in the western U.S. range, with the exception of the total lead concentration for TH-26, which exceeds the range of 2-300 mg/Kg (Bowen, 1979). Table 5 in Appendix B summarizes analytical results for total arsenic and total lead.

## 1 CONCLUSIONS AND RECOMMENDATIONS

The findings of this investigation indicate that ground water containing PCE is entering the CDOT project area near I-70 and Brighton Boulevard from the southeast, and flowing under the project area toward the northwest. The plume of PCE contamination appears to extend

from at least as far south as 44<sup>th</sup> Street and Brighton Boulevard, to at least as far northeast as the 4600 block of Brighton Boulevard, north of I-70 and Brighton Boulevard.

Ground water was again determined to flow to the northwest, confirming earlier findings. The concentrations of PCE were highest between TH-24 (530 µg/L and 480 µg/L in 1998 and 1999, respectively), and TH-26 (160 µg/L in 1999). PCE concentrations at TH-25 (23 µg/L in 1999) and at a number of locations near Brighton Boulevard and 44<sup>th</sup> Street suggest that concentrations may decrease to the northeast and southwest of the areas of highest observed PCE concentrations.

The findings of this investigation indicate ground water containing PCE at concentrations exceeding Colorado Ground Water Standards is entering the CDOT project area from upgradient sources, and that the PCE plume has extended under numerous properties. The full extent and origin of the PCE plume have not been delineated.

VALSH recommends that CDOT transmit the findings of this investigation to the Colorado Department of Public Health and Environment (CDPHE). CDPHE or perhaps the U.S. EPA may wish to conduct further investigations to determine the origin and extent of the PCE contaminated ground water in this area. The PCE contaminated ground water appears to originate from sources outside of CDOT's project area.

VALSH recommends that CDOT consider that any ground water which may be encountered during construction in the vicinity of I-70 and Brighton Boulevard may contain PCE in excess of Colorado Ground Water Standards. Any construction activities that involve handling of ground water will require prior testing of ground water for PCE, and proper handling and possibly treatment of PCE contaminated ground water.

No significant contamination by VOCs (including PCE) was detected in soil samples from 0.3 to 3.7 meters below ground surface at TH-26, TH-27 and TH-28. There was also no significant BTEX, TEPH or SVOC contamination in these soils.

Arsenic and lead concentrations were measured in surface soil (0 to 0.3 meters) at TH-26, TH-27 and TH-28. These surface soil sample locations are not at residential properties. The highest observed arsenic concentration (38 mg/Kg at TH-26) was below target levels of the U.S. EPA investigation of residential soils in this area. The highest observed lead concentration (850 mg/Kg at TH-26) exceeded the target concentration of the U.S. EPA investigation of residential soils in this area. WALSH recommends that CDOT transmit the arsenic and lead data for these surface soils to CDPHE and the U.S. EPA.

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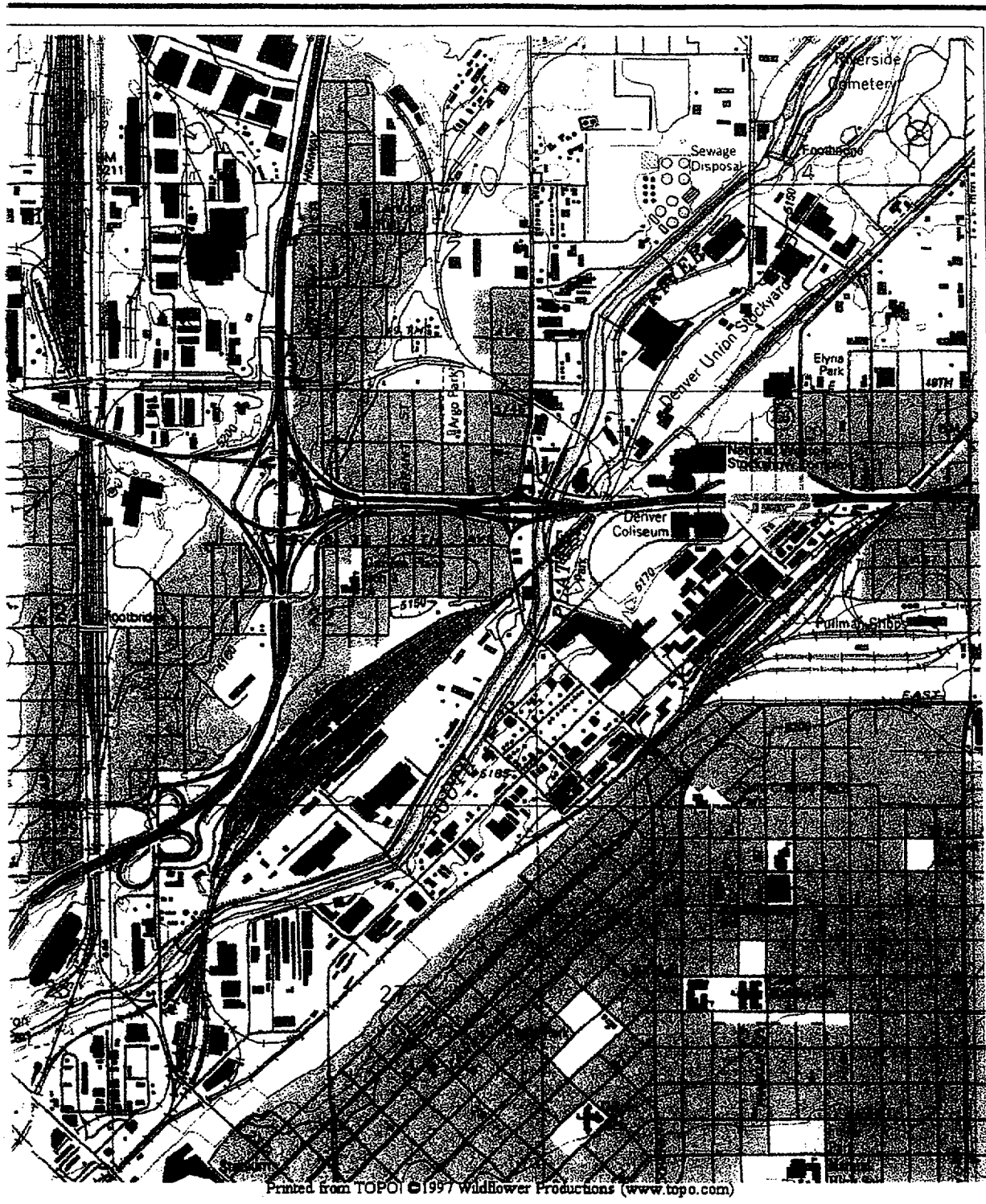
Walsh Environmental Scientists and Engineers, Inc. (WALSH), 1998. Site Investigation I-70 Phase II and III Construction 44<sup>th</sup> Street to Brighton Boulevard, *City and County of Denver, Colorado* Prepared for the CDOT, June, 1998.

Walsh Environmental Scientists and Engineers, Inc. (WALSH), 1999. Corrective Action Plan and Final Site Characterization Report Parcel 49, Tank 2, I-70 and Brighton Boulevard, Denver, Colorado, July 19, 1999.



## **APPENDIX A**

### **FIGURES**



Approximate Map Scale: 1: 15,000

R 68 W

Map Source: USGS Commerce Quadrangle  
Colorado, 7.5 Min. Series (topographic) 1959  
Revised 1980

Site



**Walsh**

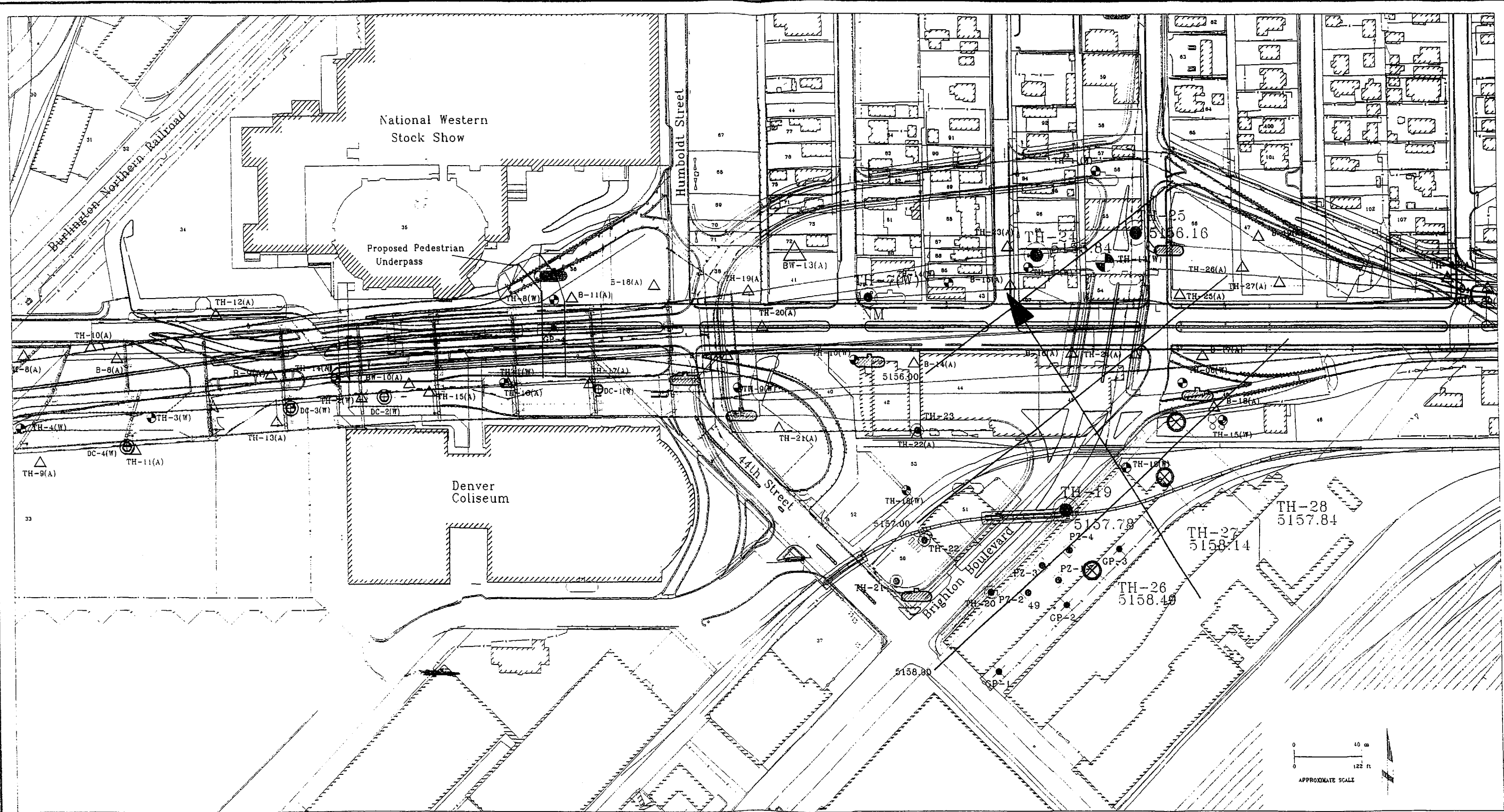
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### Site Vicinity Topographical Map

Job 3026-010

Date 9/99

Figure 1



# Explanation

- |  |  |  |  |   |
|--|--|--|--|---|
| <ul style="list-style-type: none"> <li>Washington Street Shallow Soil Boring (Walsh, 1997)</li> <li>Denver Coliseum Area Test Holes (Walsh, 1997)</li> <li>Denver Coliseum Area Test Holes (Walsh, 1997) Completed as Monitor Wells</li> <li>Walsh Monitor Well (1991)</li> <li>Aguirre Geotechnical Hole (1994, 1995)</li> <li>Walsh Geoprobe Soil Sample Site, 1998</li> </ul> | <ul style="list-style-type: none"> <li>Approximate Property Boundary and ID Number</li> <li>Former Aboveground Storage Tank(s)</li> <li>Walsh Test Hole, 1999 Potentiometric Surface Measured on 8/25/99</li> <li>Former Barrel Storage</li> </ul> | <ul style="list-style-type: none"> <li>Former Underground Storage Tank(s)</li> <li>Underground Storage Tank(s) - Removal Unknown</li> <li>Underground Storage Tank(s)</li> <li>Walsh Temporary Piezometer, 1999 Sampling Location</li> <li>Potentiometric Surface Contour Contour Interval = 1.0 foot</li> </ul> | <ul style="list-style-type: none"> <li>Proposed Monitoring Well</li> <li>Walsh Test Hole, 1998, 1999 Resample Location, Potentiometric Surface Measured on 3/23/99</li> <li>Aguirre Geotechnical Hole (1994, 1995) Sampling Location</li> <li>Ground Water Flow Direction</li> </ul> | <ul style="list-style-type: none"> <li>Proposed sidewalks and riprap</li> <li>Proposed curbs and gutters</li> <li>Proposed retaining walls</li> <li>Proposed storm and/or sanitary sewers</li> <li>Proposed gas</li> <li>Proposed telephone</li> <li>Proposed water line</li> </ul> |
|--|--|--|--|---|

**Walsh**  
Environmental Scientists and Engineers, Inc.  
Potentiometric Surface Map  
Test Hole Locations

Job 3026-010 | Date 6/99 | Figure 4

## **APPENDIX B**

### **TABLES**

**Table 1 - Ground Water Elevation Data**

Monitoring Well	North Ground Level Elevation (Ft.)	North TOC <sup>2</sup> (Ft.)	Ground Water Level* (Ft. below TOC)	Potentiometric Surface Elevation (Ft.)
TH-7	NM	NM	27.80*	-
TH-19	5187.58	5187.27	29.49*	5157.78
TH-24	5185.21	5185.11	29.27*	5155.84
TH-25	5185.14	5184.98	28.82*	5156.16
TH-26	5186.62	5190.19	31.70**	5158.49
TH-27	5188.14	5191.74	33.60**	5158.14
TH-28	5186.10	5189.64	31.80**	5157.84

No Measurement

Top of Casing

Ground water levels measured on 03/23/99

Ground water levels measured on 06/25/99

**Table 2 - PID Soil Screening Measurements**

Monitoring Well	Date	Depth Meters (feet. bgs.) <sup>1</sup>	PID Reading (ppm) <sup>2</sup>
TH-26	06/16/99	0-0.15 (0-5)	0.3
	06/16/99	1.5-2.1 (5-7)	2.4
	06/16/99	3-3.7 (10-12)	7.0
	06/16/99	4.6-5.2 (15-17)	0.0
	06/16/99	6.1-6.7 (20-22)	0.0
	06/17/99	7.6-8.2 (25-27)	0.0
	06/17/99	9.1-9.8 (30-32)	0.0
	06/18/99	0-0.3 (0-1)	12.8
TH-27		1.5-2.1 (5-7)	23.6
		3-3.7 (10-12)	6.9
		4.6-5.2 (15-17)	5.5
		6.1-6.7 (20-22)	1.5
		7.6-8.2 (25-27)	0.9
		9.1-9.8 (30-32)	0.1
	06/17/99	0-0.3 (0-1)	0.0
		0.3-0.9 (1-3)	1.4
TH-28		1.5-2.1 (5-7)	1.6
		3-3.7 (10-12)	3.9
		4.6-5.2 (15-17)	0.6
		6.1-6.7 (20-22)	0.0
		7.6-7.8 (25-27)	0.0
		9.1-9.8 (30-32)	0.0

**Table 3 – VOCs and TEPH Soil Analytical Results (mg/Kg)**

Monitoring Well	Date	VOCs	VOCs	TEPH	TEPH
TH-26	6/16 & 6/17/99	ND @ 0.3-0.9 (1-3) ##	ND @ 7.6-8.2 (25-27)	11 @ 0-0.3 (0-1)	ND @ 7.6-8.2 (25-27)
TH-27	6/18/99	ND @ 1.5-2.1 (5-7)	ND @ 7.6-8.2 (25-27)	ND @ 1.5-2.1 (5-7)	ND @ 7.6 8.2 (25-27)
TH-28	6/17/99	ND @ 3.0-3.7 (10-12)	ND @ 7.6-8.2 (25-27)	6 @ 3.0-3.7 (10-12)	ND @ 7.6-8.2 (25-27)

- VOCs include BTEX analyses  
# Depth in Meters bgs (feet bgs)

**Table 4 – Semi-Volatile Organic Compounds (SVOCs) Soil Analytical Results (mg/Kg)**

Monitoring Well	Date	Phenathrene	Fluoranthene	Pyrene	Chrysene	Benzo(b)-Fluoranthene
TH-26 @ 0-0.3 (0-1) ##	6/16 & 6/17/99	<0.33	<0.33	<0.33	<0.33	<0.33
TH-27 @ 0-0.3 (0-1)	6/18/99	<0.33	<0.33	<0.33	<0.33	<0.33
TH-28 @ 0.3-0.9 (1-3)	6/17/99	0.34	0.56	0.92	0.37	0.43

# = Depth in meters (feet bgs)

**Table 5 – Total Arsenic and Lead Soil Analytical Results (mg/Kg)**

Monitoring Well	Date	Total Arsenic	Total Lead
TH-26 @ 0-0.3 (0-1) ##	6/16/99	38	850
TH-27 @ 0-0.3 (0-1)	6/18/99	4.7	68
TH-28 @ 0.3-0.9 (1-3)	6/17/99	15	140
Western U.S. Range*	1979	0.1-40	2-300

Concentrations in BOLD exceed Western U.S. Range background concentrations  
\* Bowen 1979

**sh Environmental, Inc.**  
**LOG**

Location Sketch or Description

OLD #2

TH-27

# STAKE  
04124

4N

Boring Number  
TH-27

Sheet

1 of 1

Drilling Contractor  
CHS SS

Location 162, 27 Sub # 424124  
Drilling Contractor SITE SVCS 0 600FF / TONY

Start Date 6/18/99

Finish 6/18/99

Logger V.A. S. 000000

Sample			Standard Penetration Test Results	Soil Description	Symbol of USCS Log	Comments
Interval	Tag No.	Recovery				
			6" - 6" - 6"	USCS Group Symbol, Name, Gradation or Plasticity, Particle Size Distribution, Color, Moisture Content, Relative Density or Consistency, Soil Structure, Mineralogy.		PID readings/stainings
			(2)			
			0-1 2809	Grey-BK Silty clay, w/ Sand, F. No odors or stains	CL	PID = 12.8 (SOILS / METER)
			20818 3-4-4-4 24/24"	Similar a.g. to 5'3" Into Sand = SILT, F, grading to Sand, F, grey/orange @ 6.5'. Loose, damp. No odors/stains	SM /SA	PID = 23.6 (SOILS / METER) (FTH) (CL) in shear (grey)
			20824 7-18-20-21 24/24"	lt br. Sand c-m; some F, gravel, subdr. F, loose, damp. No odors or stains	SP	PID = 6.9
			20832 12-28-31-31 24/24"	Br Silt/Sand, Fine w/ F gravel to 15'6" Into c-vc sand, to m-F sand w/ cobbles & F gravel, loose, dry.	SM /SP	PID = 5.5
			20844 2-18-21-22-21 16/24"	a.g. (15.6 - 17')	SA	PID = 1.5
			20854 12-28-22-21 /24"	lt br-wt Sand, m-F, fine gravel, & cobbles coarsening towards 27'. Damp to moist (Fringe). NO odors/stains	SA	PID = 0.9 V 225-27' = (Fringe) (NOILS) (FTH)
			20906 4-16-29-39 /24"	Gravel, m-F, sandy, & vc sand. Some m-F, saturated to ~31'. & Predominantly c-m Sand, br, some F, & F gravel to 32'. Saturated EOB = 30' Sit Well	GW	PID = .1

**LOG**

Boring Number **TH-28** Sheet **1** of **1**

Location Sketch or Description



Location \_\_\_\_\_  
 Drilling Contractor **SITE SICS**  
 Method and Equipment \_\_\_\_\_  
 and Date \_\_\_\_\_ Start **6/17/94** Finish \_\_\_\_\_ Logger **VAS**

Sample			Standard Penetration Test Results	Soil Description	Symbol of USCS Log	Comments
Interval	Tag No.	Recovery				
			6" - 6" - 6" (N)	USCS Group Symbol, Name, Gradation or Plasticity, Particle Size Distribution, Color, Moisture Content, Relative Density or Consistency, Soil Structure, Mineralogy.		PID readings/stainings
				0-1' - BIK SM, silt/sand, little clay, could dust	SM/CL	PID = 0.0
				1-3' a.a. but wet		PID = 1.4 (PL-As) (VOC)
				3-5' a.a. but Br, no could dust		
			@1153 2-2-2-3 24/24	as 0-1 to 5'3" into SM silty sand, with clay to 2'5" $\downarrow$ clay, & extend to 7' with $\uparrow$ fine gravel. Damp.	SM -ML	PID = 1.6
			@1204 5-6-7-8 24/24	Sand, m-F, with F gravel, br, loose, dry to 10'6" into SM silty sand, br, silty plastic, damp w/ $\uparrow$ clay towards 12'. No odor / stains	SP SM	PID = 3.9 (VOC) (TEH)
			@1215 8-8-18-20 24/24	lt. br-tan Sand, F, & silt, m-F gravel, dry, loose to 16' 16'-17' m-F gravel to m-F sand, loose, white-br, dry. No odor / stains. (2" of Fe stain)	SP SM GW	PID = 2.6
			@1223 4-9-12-17 23/24	lt br sand, m-F + fine gravel, loose, dry to 21'4" then sand, with 7 m-F gravel & cobbles. 1" of clay @ ~22'. No stain or odor.	SP GW	PID =
			@1236 6-12-15-21 19/24"	Sand, lt br-white, m-F, some c sand, m-F gravel & cobbles. LOOSE. MOIST. No odor / stains	SP	PID ~25' (VOC) (TEH)
			@1245 24/24 12-15-16-14	A.A. but saturated.  Set well @ 30'		



## WELL CONSTRUCTION LOG

.3026-010 Well No. TH-26

### Drilling Summary

Total Depth of Hole: 30'  
 Hole Diameter: 7" OD  
 Drilling Company: SITE SVCS  
 Driller: TONY / IGOTT  
 Rig Type: CME-55  
 Bits: 7" HSA, 2' LONG SS SAMPLE  
 Geologist: KA. SECORNO

### Construction Time Log

	Start		Finish	
	Date	Time	Date	Time
Drilling:	_____	_____	_____	_____
Screen Placement:	_____	_____	_____	_____
Filter Placement:	_____	_____	_____	_____
Seal Placement:	_____	_____	_____	_____
Grouting:	_____	_____	_____	_____

### Depth to Water

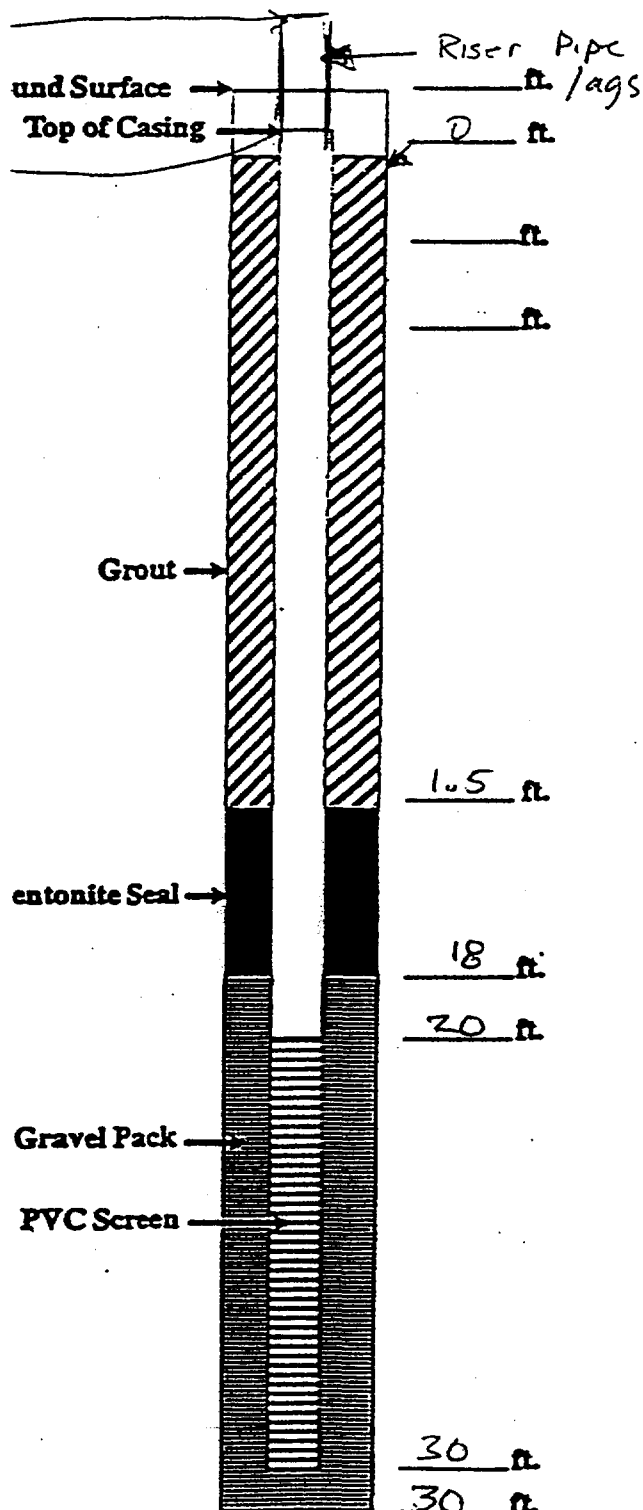
Depth: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

### Well Construction Materials

	Grout	Seals	Filter
Quantity:	_____	_____	_____
Type:	_____	_____	_____
Screen			
Size:	<u>10</u>	Config:	<u>.010 5/62</u>
Area/Ft.	<u>17 gal/ft</u>	Comp:	<u>sch 40</u>
Inside Diam:	<u>2"</u>	Outside Diam:	<u>PVC</u>

### Comments

Riser Protective M/O unit



Measuring point is top of CASING  
 unless otherwise noted

## WELL CONSTRUCTION LOG

Job No. 3026-010 Well No. TH-27

### Drilling Summary

Total Depth of Hole: 30'  
Hole Diameter: \_\_\_\_\_  
Drilling Company: SITE SERVICES  
Driller: TONY / GCOFF  
Rig Type: CME-55  
Bits: HSA, 2' SS samplers  
Geologist: V. D. SECONAZ

### Construction Time Log

	Start		Finish	
	Date	Time	Date	Time
Drilling:	_____	_____	_____	_____
Screen Placement:	_____	_____	_____	_____
Filter Placement:	_____	_____	_____	_____
Seal Placement:	_____	_____	_____	_____
Grouting:	_____	_____	_____	_____

### Depth to Water

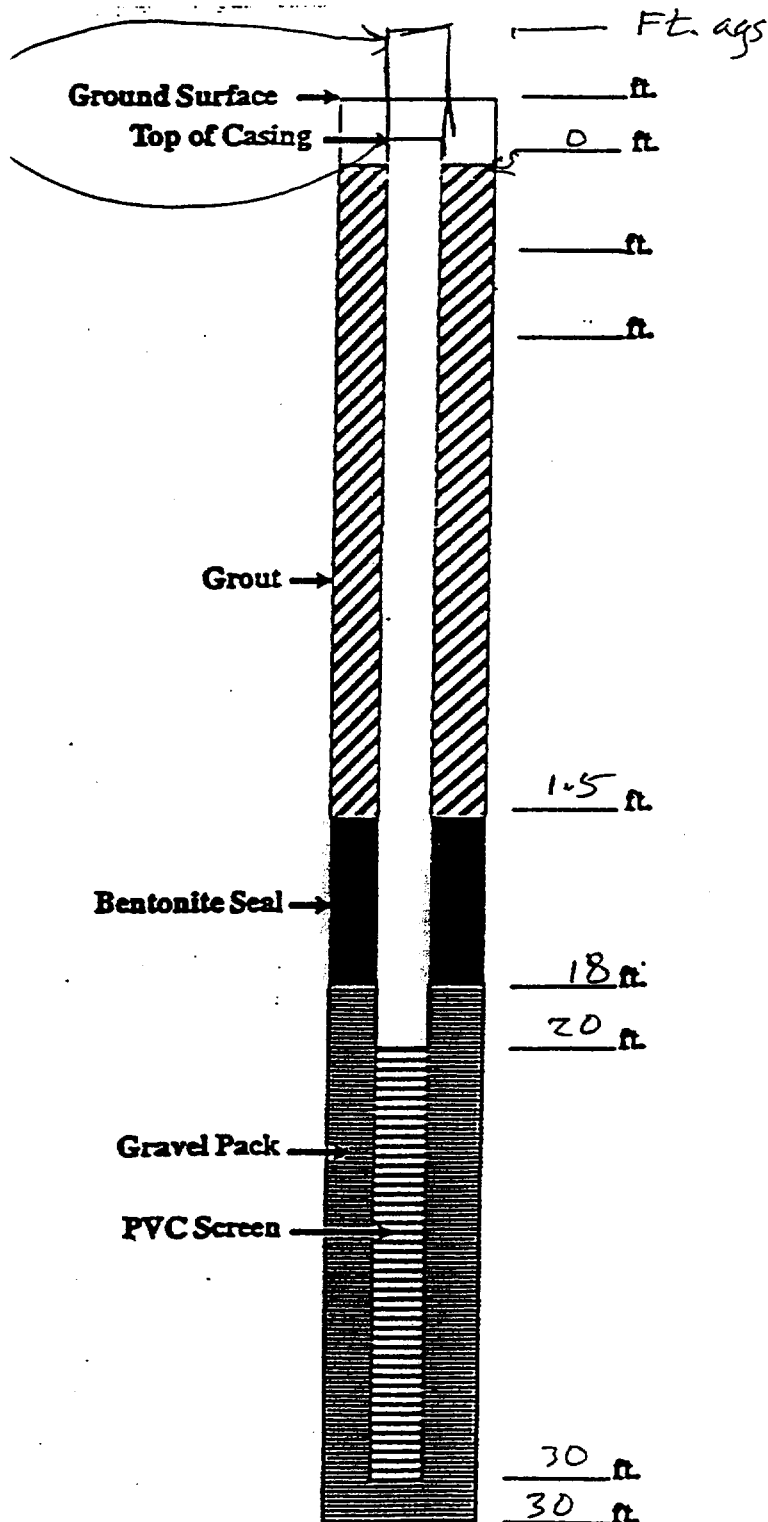
Depth: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

### Well Construction Materials

	Grout	Seals	Filter
Quantity:	_____	_____	_____
Type:	_____	_____	_____
Screen			
Size:	<u>10'</u>	Config:	<u>.010 slot</u>
Area/Ft:	<u>17 gal/Ft</u>	Comp:	<u>sch 40</u>
Inside Diam:	<u>2"</u>	Outside Diam:	<u>PVC</u>

### Comments

Riser Protective Mount



Measuring point is top of CASING  
unless otherwise noted

## WELL CONSTRUCTION LOG

3026-C10 Well No. TH-28

### Drilling Summary

Total Depth of Hole: 30'  
 Hole Diameter: \_\_\_\_\_  
 Drilling Company: SITE SERVICES  
 Driller: JUNY / GCOFF  
 Rig Type: CMS-55  
 Bits: HSA, 2' SS sampler  
 Geologist: V. A. SECORADO

### Construction Time Log

	Start		Finish	
	Date	Time	Date	Time
Drilling:	_____	_____	_____	_____
Screen Placement:	_____	_____	_____	_____
Filter Placement:	_____	_____	_____	_____
Seal Placement:	_____	_____	_____	_____
Grouting:	_____	_____	_____	_____

### Depth to Water

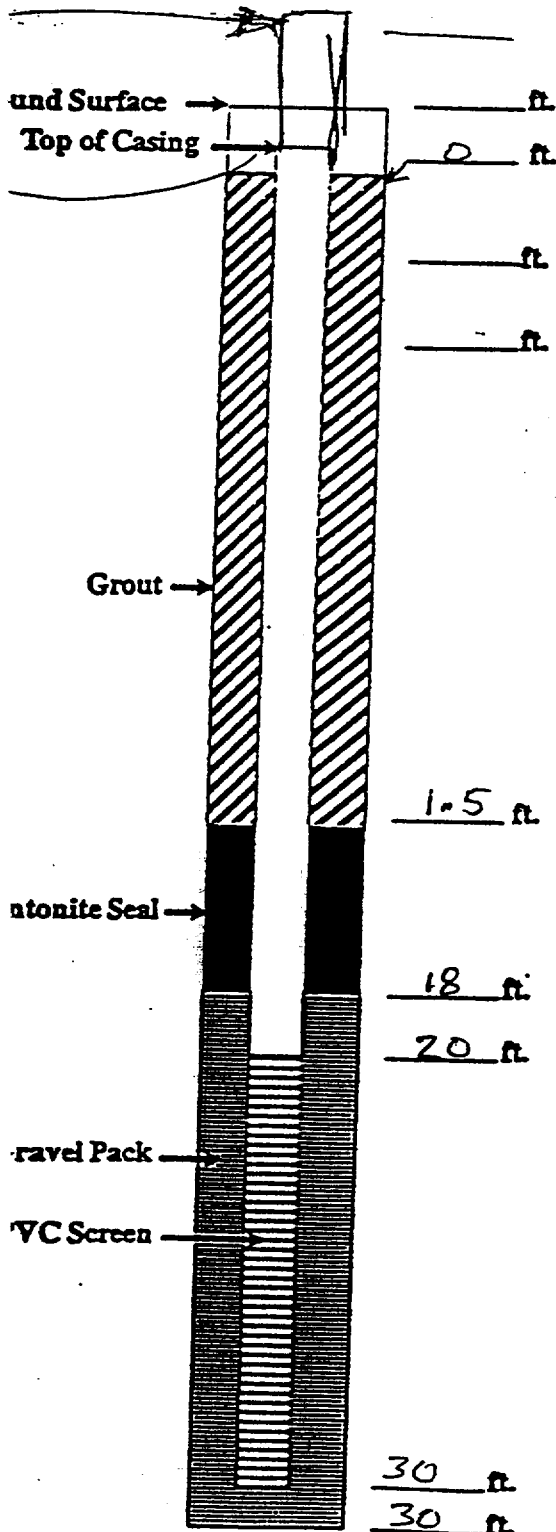
Depth: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

### Well Construction Materials

	Grout	Seals	Filter
Quantity:	_____	_____	_____
Type:	_____	_____	_____
Screen			
Size:	<u>10'</u>	Config:	<u>0.010 slot</u>
Area/Ft	<u>17.9 sq/Ft</u>	Comp:	<u>Sch 40</u>
Inside Diam:	<u>2"</u>	Outside Diam:	<u>AW</u>

### Comments

Riser Protective Mount



Measuring point is top of CASING  
 unless otherwise noted

## **APPENDIX D**

### **FIELD DATA TRANSMITTAL SHEETS**

# GROUND-WATER SAMPLING FIELD DATA SHEET

Well Number TH-26  
 Well Diameter 2"  
 Well Vals  
 Well Number 58378, 79, 80 VAS  
 58390, 91, 92

Project Number 3026-010  
 Project Name I-70 / Brighton  
 Date 6/23/99  
 Time 0800

Well Number	Sampling kit	pH 7.00			
46"	Hydac #3	Actual	Adjust	Temp	Time
	Solinst #2				
	MMC				
Water Level (ft)	Conductivity	pH 10.00			
31.7	Std =				
	Actual = @ °F				
		Actual	Adjust	Temp	Time
Depth (ft)	Clear Bailer Result	Purging Equipment			
34.3		DISP. Bailer			
Thickness	Sample Depth (ft)	Sampling Equipment			
2.6		" "			
Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
.4					

Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
0	INITIAL	67	1060	6.9	Silty Brackish no odor
1	.4	65.8	940	6.11	Silty Brackish no odor
2	.8	63.1	920	5.88	Silty Brackish no odor
3	1.2	62.6	900	5.72	Silty Brackish no odor
4	1.6	61.5	880	5.64	Silty Brackish no odor
5	2.0	62.2	960	5.59	Silty Brackish no odor
↓					
Sample					

3:

Collected by: VAS/EC  
 Checked by:

for:

W/TEH	BTEX	THE	TVH	OTHER
		X		VAS

# GROUND-WATER SAMPLING FIELD DATA SHEET

Number TH-27

1g Diameter 2"

onnel VPS

Number 58393, 58394, 58395

Project Number 3026-010

Project Name I-70/Brighton

Date 2/23/99

Time 0830

Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
46"	Hydac #3				
	Solinst #2				
	MMC				
Water Level (ft)	Conductivity Time	pH 10.00			
33.6	Std =				
	Actual = @ °F				
		Actual	Adjust	Temp	Time
Well Depth (ft)	Clear Bailer Result	Purging Equipment			
34.1		DISP. Bailer			
Thickness	Sample Depth (ft)	Sampling Equipment			
.5		" "			
Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				

Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
0					
1					
2					
3					
4					
5					
↓					
Sample					

Notes: NO C.V.'S REMOVED. Total samples due to low H<sub>2</sub>O quantity.

Collected by: VPS/EC

Checked by:

1 for: TEH Sample 3/4 Full

MT/TEH	BTEX	THE	TVH	OTHER
		X		VOCs

# GROUND-WATER SAMPLING FIELD DATA SHEET

Number TH-28

ing Diameter 2"

onnel VPS

Number 54684, 85, 86 VPS

58396, 97, 98

Project Number 3026-010

Project Name I-70/Brighton

Date 6/23/99

Time 0914

Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
	Hydac #3				
	Solinst #2				
	MMC				
Water Level tickup) (ft)	Conductivity Time Std = Actual = @ °F	pH 10.00			
31.8		Actual	Adjust	Temp	Time
Well Depth ckup) (ft)	Clear Bailer Result	Purging Equipment			
34.08		DPA. Bailer			
d Thickness	Sample Depth (ft)	Sampling Equipment			
2.28		" "			
Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
4					

Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
0	Initial	70.2	1240	5.93	clear slightly turbid no odor
1	.4	65.4	1140	5.57	Brown silty no odor
2	.8	63.6	1120	5.49	Brown/silty no odor
3	1.2	63.2	1090	5.39	Brown/silty no odor
4	1.6	62.4	1100	5.33	Brown/silty no odor
5	2.0	63.4	1060	5.33	Brown silty no odor
↓					
Sample					

(S:

Collected by: VPS

Checked by:

d for:

VH/TEH	BTEX	THE	TVH	OTHER
		X		BZ60 VOCs

# GROUND-WATER SAMPLING FIELD DATA SHEET

Lumber ~~TH-7(W)~~  
 g Diameter 2"  
 nnel VAS  
 umber

Project Number 3026 010  
 Project Name I-70 / Brighton Blvd.  
 Date 3/23/99  
 Time

Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
	Hydac #2				
	Solinst #5				
	MMC				
Water Level (ft)	Conductivity	pH 10.00			
27.80	Std =				
	Actual = @ °F				
		Actual	Adjust	Temp	Time
Well Depth (ft)	Clear Bailer Result	Purging Equipment			
37.91		Disp. Bailer			
Well Thickness	Sample Depth (ft)	Sampling Equipment			
10.11		" "			
Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
1.7					

Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
INITIAL	0	62.7	8960	11.15	Clear
1	1.7	58.1	8780	8.79	" color??
2	3.4	56.0	7820	8.75	Turbid, Lt. br color??
3	5.1	56.1	7380	8.64	" "
Sample					

KS:

Collected by: VAS  
 Checked by:

Send for:

TVH/TEH	BTEX	THE	TVH	OTHER
		X	X	8260 VOCs



# GROUND-WATER SAMPLING FIELD DATA SHEET

Well Number **TH-19**  
 Casing Diameter **2" ID**  
 Personnel **VAS**  
 Number

Project Number **3026-CID**  
 Project Name **I-704 Brighton Blvd.**  
 Date **03/23/99**  
 Time **0734**

g Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
	Hydac #2	7.81	7.0	38.2	0740
	Solinst #5				
	MMC				
Water Level (stickup) (ft)	Conductivity Time	pH 10.00			
<b>29.49</b>	Std = 100.5				
	Actual = 1010 @ 40.0 PF				
		Actual	Adjust	Temp	Time
		10.21	10.0	39.1	0741
Well Depth (stickup) (ft)	Clear Bailer Result	Purging Equipment			
<b>33.9</b>		<b>Disp. Bailer</b>			
ated Thickness	Sample Depth (ft)	Sampling Equipment			
<b>4.41</b>		" "			
g Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
<b>.75</b>					

ne	Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
58	INITIAL	0	52.3	1490	9.91	Clear, site turning towards tip. NO H.C. color.
00	1	.75	54.0	1490	9.21	Rust orange NO H.C. color.
02	2	1.5	54.7	1550	9.11	" "
06	3	2.25	54.8	1620	9.09	" "
	Sample					

arks:

Collected by: **VAS**  
 Checked by:

pled for:

EX/TVH/TEH	BTEX	THE	TVH	OTHER
		X	X	8260 VOCs

# GROUND-WATER SAMPLING FIELD DATA SHEET

Number *TH-24*  
 ng Diameter *2" ID*  
 onnel *VPS*  
 Number

Project Number *3026-010*  
 Project Name *I-70 / Brighton Blvd.*  
 Date *2/23/99*  
 Time *0955*

Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
	Hydac #2				
	Solinst #5				
	MMC				
Water Level (ft)	Conductivity Time	pH 10.00			
<i>29.27</i>	Std =				
	Actual = @ °F				
		Actual	Adjust	Temp	Time
Well Depth (ft)	Clear Bailer Result	Purging Equipment			
<i>35.10</i>		<i>Disp. Bailer</i>			
sed Thickness	Sample Depth (ft)	Sampling Equipment			
<i>5.83</i>		<i>" "</i>			
Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
<i>1.0</i>					

	Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
5	INITIAL	0	64.6	1920	9.61	<i>clear to set. finished, NO. 1000 Rustorange.</i>
7	1	1	64.5	1810	9.47	
2	2	2	60.9	1950	9.47	
3	3	3	61.0	1840	9.37	<i>" "</i>
5	Sample					

rks:

Collected by: *VPS*  
 Checked by:

led for:

QTVH/TEH	BTEX	THE	TVH	OTHER
		<i>X</i>	<i>X</i>	<i>8260 VOCs</i>

# GROUND-WATER SAMPLING FIELD DATA SHEET

Well Number **TH-25**  
 Casing Diameter **2" I.D.**  
 Personnel **VAS**  
 Log Number

Project Number **3026-010**  
 Project Name **I-70 / Brighton Blvd**  
 Date **3/23/99**  
 Time **1644**

Initial Stickup	Sampling kit	pH 7.00			
		Actual	Adjust	Temp	Time
	Hydac #2				
	Solinst #5				
	MMC				
Static Water Level (ft)	Conductivity Time	pH 10.00			
<b>28.82</b>	Std =				
	Actual = @ °F				
		Actual	Adjust	Temp	Time
Well Depth (ft)	Clear Bailer Result	Purging Equipment			
<b>34.90</b>		<b>Dist. Bailer</b>			
Gravel Thickness	Sample Depth (ft)	Sampling Equipment			
<b>6.16</b>		" "			
Log Volume	REMEMBER: 2" well multiply by .17 4" well multiply by .66				
<b>1.0</b>					

Time	Casing Volumes	Gallons Removed	Temperature °F	Conductivity uS/cm	pH	Comments
54	INITIAL	0	63.5	1960	9.25	clear
58	1	1.0	58.3	1810	9.50	Rust-orange
2	2	2.0	58.1	1870	9.10	" "
65	3	3.0	57.5	1840	9.07	" "
77	Sample					

Remarks:

Collected by: **VAS**  
 Checked by:

Sampled for:

TEX/TVH/TEH	BTEX	THE	TVH	OTHER
		X	X	8260 VOCs

## **APPENDIX E**

### **SURVEY DATA**

C7419 Walsh Environmental Scientists & Engineers (Proj. No. 3026-010)

Unit: fts  
Coordinate Type: Geodetic  
Reference Ellipsoid: WGS-1984  
Projection Set: CO-CEN-83  
Datum: NAD-83 ('92 Adj)/NAVD-88

#301, 39-46-46.08393 N, 104-57-59.67069, 5189.638 Top of PVC  
TH-28  
#351, 39-46-43.08319 N, 104-57-59.66592, 5186.095  
Set "X" on Concrete Base

#302, 39-46-45.01935 N, 104-58-00.26886, 5191.736 Top of PVC  
TH-27  
#352, 39-46-45.02431 N, 104-58-00.27362, 5188.142  
Set "X" on Concrete Base

#303, 39-46-43.45441 N, 104-58-01.53681, 5190.188 Top of PVC  
TH-26  
#353, 39-46-43.45237 N, 104-58-01.53782, 5186.619  
Set "X" on Concrete Base

#311, 39-46-42.47983 N, 104-58-03.96050, 5189.842 Top of PVC  
PZ-1

#312, 39-46-42.47823 N, 104-58-03.23252, 5190.608 Top of PVC  
PZ-2

#313, 39-46-42.81011 N, 104-58-03.64789, 5190.084 Top of PVC  
PZ-3

#314, 39-46-42.98182 N, 104-58-03.26159, 5190.440 Top of PVC  
PZ-4



company

Environmental Inc.  
Pearl East Circle #108  
er CO 80301

Stan Spencer

325 Interlocken Parkway  
Suite 200  
Broomfield, CO 80021  
(303) 469-8868  
(800) 873-8707  
FAX: (303) 469-5254

Order #: 99-06-215  
Date: 06/27/99 16:15  
Work ID: 3026-010 44TH & BRIGHTON  
Date Received: 06/17/99  
Date Completed: 06/24/99

#### SAMPLE IDENTIFICATION

<u>Client Description</u>	<u>Sample</u>	<u>Client Description</u>
	<u>Number</u>	
TH-26 TAG #58380	02	TH-28 TAG #58382

ed are the analytical results for the submitted sample(s). Please  
the CASE NARRATIVE for a discussion of any data and/or quality  
l issues. A listing of data qualifiers and analytical codes is  
d on the TEST METHODOLOGIES page at the end of the report.

have any questions regarding the analyses, please feel free to call.

ely,

K. Toon  
Manager

es were prepared and analyzed according to methods outlined in the  
wing references:

Test Methods for Evaluating Solid Waste, USEPA SW-846, Third Edition,  
Revision 4, December 1996.

Standard Method for Laboratory Determination of Water (Moisture) Content of  
Soil, Rock, and Soil-Aggregate Mixtures, ASTM D 2216-80, July 1980.

ems encountered with the analyses are discussed in the following narrative.

atrix spike recovery for the batched lead quality control sample was below  
ower control limit. All other quality assurance analyses were acceptable.

## **APPENDIX F**

### **LABORATORY DATA & CHAIN OF CUSTODY**



Sample: 01A TH-26 TAG #58380

Collected: 06/16/99

Matrix: SOIL

<u>Test Description</u>	<u>Method</u>	<u>Result</u>	<u>Q</u>	<u>Limit</u>	<u>Units</u>	<u>Analyzed</u>
Arsenic, Total	SW 7060	38	DS	2.4	mg/Kg-DRY	06/23/99
Lead, Total	SW 7421	850	D	24	mg/Kg-DRY	06/23/99
Percent Moisture	ASTM D2216	16.5		0.1	WT%	06/23/99

Sample: 02A TH-28 TAG #58382

Collected: 06/17/99

Matrix: SOIL

<u>Test Description</u>	<u>Method</u>	<u>Result</u>	<u>Q</u>	<u>Limit</u>	<u>Units</u>	<u>Analyzed</u>
Arsenic, Total	SW 7060	15	DS	1.4	mg/Kg-DRY	06/23/99
Lead, Total	SW 7421	140	D	2.9	mg/Kg-DRY	06/23/99
Percent Moisture	ASTM D2216	30.6		0.1	WT%	06/23/99

THE FOLLOWING CODES APPLY TO THE ANALYTICAL REPORT

RESULT field...

ND = not detected at the reported limit

NA = analyte not applicable (see case narrative/methods for discussion)

(qualifier) field...

GENERAL:

\* = Recovery or %RPD outside method specifications

H = value is estimated due to analysis run outside EPA holding times

E = reported concentration is above the instrument calibration range

D = analyte was diluted to bring within instrument calibration range or  
to remove matrix interferences

ORGANIC ANALYSIS DATA QUALIFIERS:

B = analyte was detected in the laboratory method blank

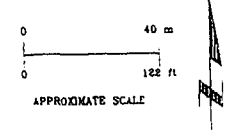
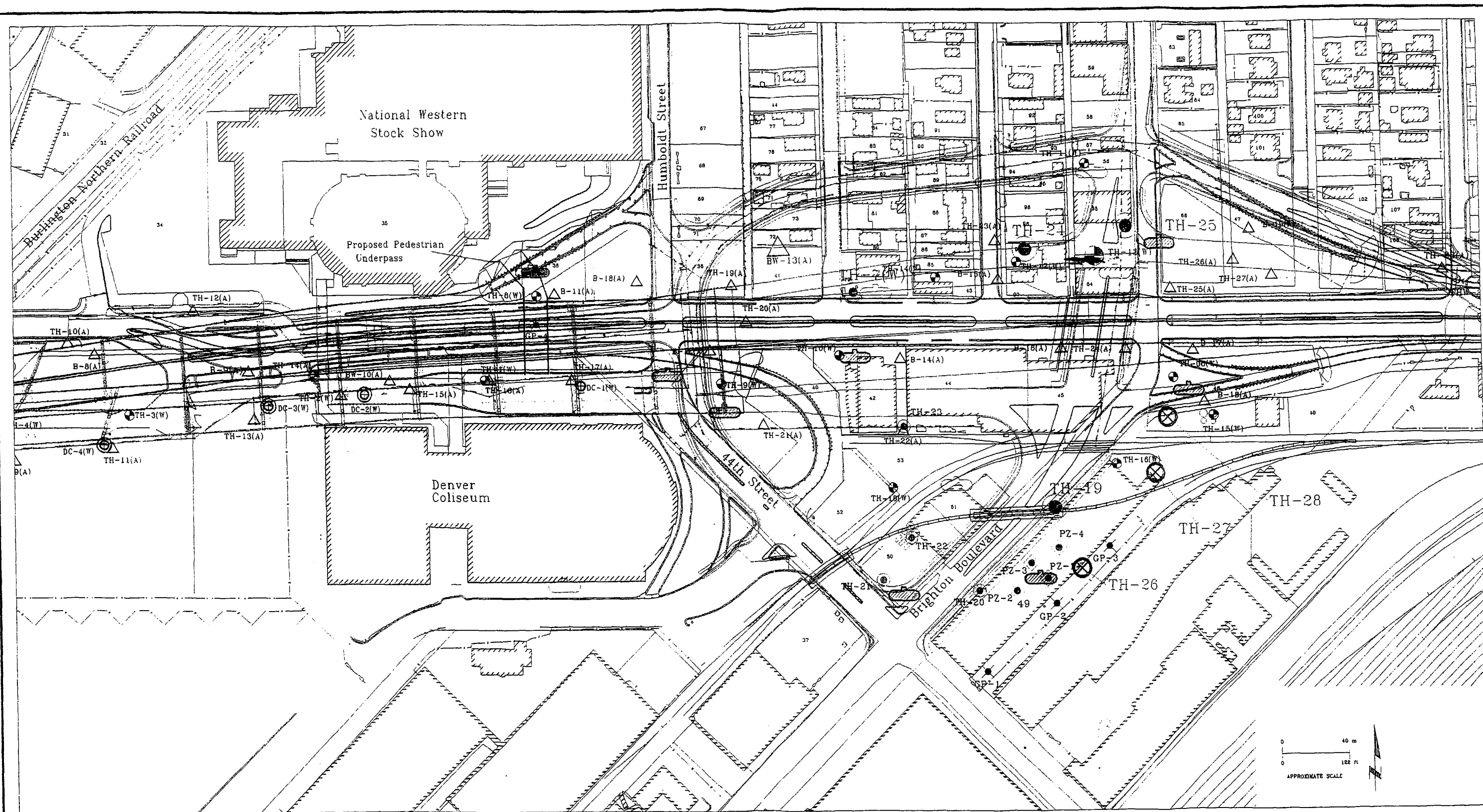
J = analyte was detected above the instrument detection limit (IDL)  
but below the analytical reporting limit (CRDL)

INORGANIC ANALYSIS DATA QUALIFIERS:

B = analyte was detected above the instrument detection limit (IDL)  
but below the analytical reporting limit (CRDL)

W = post digestion spike did not meet criteria (85-115%)

S = reported value determined by the Method of Standard Additions



# Explanation

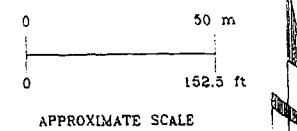
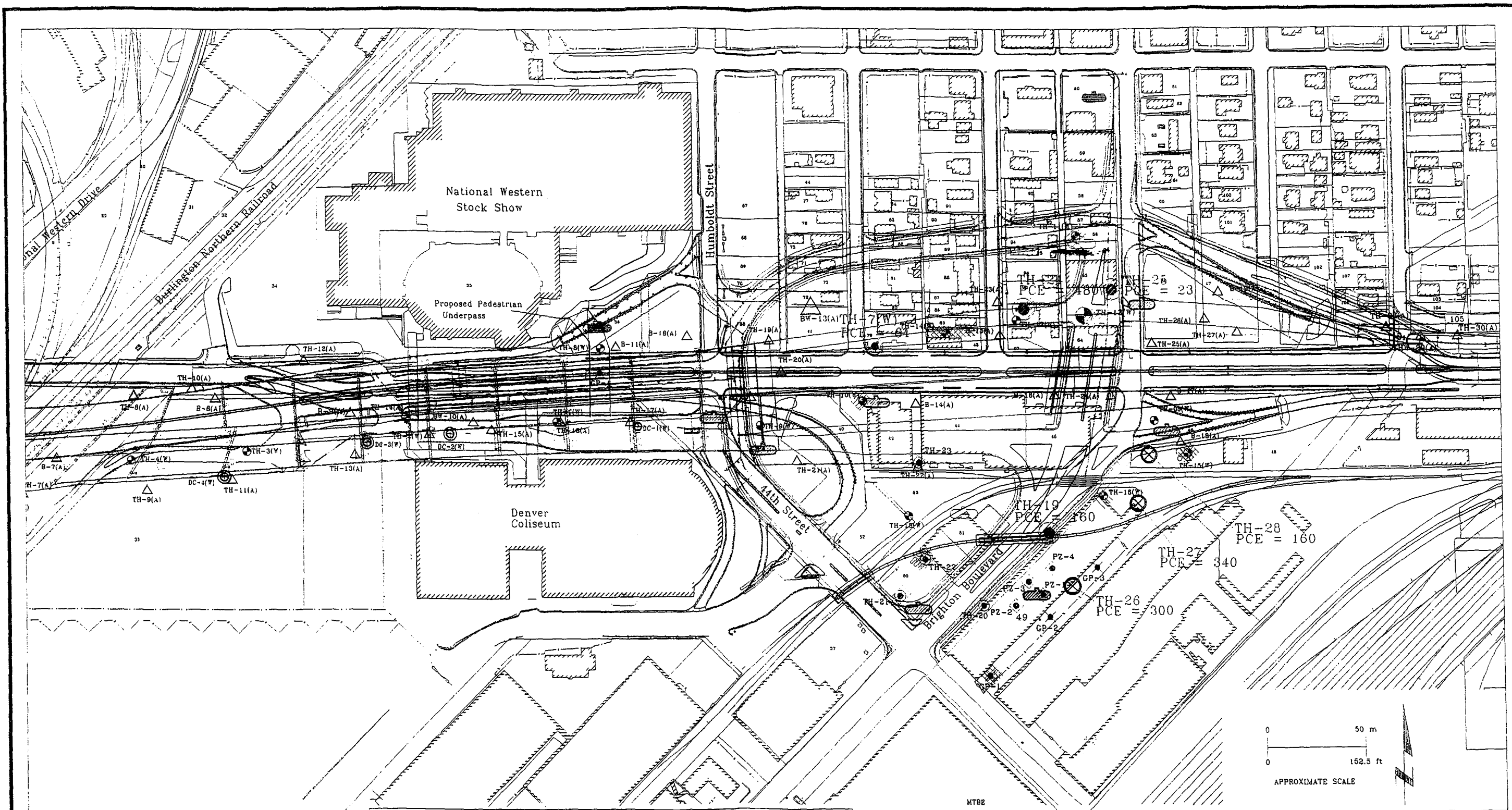
- |   |   |   |   |   |
|---|---|---|---|---|
| <ul style="list-style-type: none"> <li>Washington Street Shallow Soil Boring (Walsh, 1997)</li> <li>Denver Coliseum Area Test Holes (Walsh, 1997)</li> <li>Denver Coliseum Area Test Holes (Walsh, 1997) Completed as Monitor Wells</li> <li>Walsh Monitor Well (1991)</li> <li>Aguirre Geotechnical Hole (1994, 1995)</li> </ul> | <ul style="list-style-type: none"> <li>Walsh Geoprobe Soil Sample Site, 1998</li> <li>Walsh Test Hole, 1998</li> <li>Approximate Property Boundary and ID Number</li> <li>Former Aboveground Storage Tank(s)</li> </ul> | <ul style="list-style-type: none"> <li>Former Underground Storage Tank(s)</li> <li>Underground Storage Tank(s) - Removal Unknown</li> <li>Underground Storage Tank(s)</li> <li>Former Barrel Storage</li> <li>Approximate Location of Former Gas Station</li> <li>Tank Type (UST or AST) Unknown, UST Removal Unknown</li> <li>Walsh Test Hole, 1999</li> </ul> | <ul style="list-style-type: none"> <li>Proposed Monitoring Well</li> <li>Walsh Monitoring Well (1991) Sampling Location</li> <li>Walsh Test Hole, 1998 Sampling Location</li> <li>Aguirre Geotechnical Hole (1994, 1995) Sampling Location</li> </ul> | <ul style="list-style-type: none"> <li>Proposed sidewalks and riprap</li> <li>Proposed curbs and gutters</li> <li>Proposed retaining walls</li> <li>Proposed storm and/or sanitary sewers</li> <li>Proposed gas</li> <li>Proposed telephone</li> <li>Proposed water line</li> </ul> |
|---|---|---|---|---|

**Walsh**

Environmental Scientists and Engineers, Inc.

Site Map With  
Test Hole Location

Job 3026-010 | Date 9/99 | Figure 2



# Explanation

- |   |  |  |   |  |
|---|--|--|---|--|
| <p>WSB-(W) Washington Street Shallow Soil Boring (Walsh, 1997)</p> <p>DC-(W) Denver Coliseum Area Test Holes (Walsh, 1997)</p> <p>DC-23-(W) Completed as Monitor Wells</p> <p>TH-(W) Walsh Monitor Well (1991)</p> <p>TH-(A) Aguirre Geotechnical Hole (1994, 1995)</p> | <p>CP-1 Walsh Geoprobe Soil Sample Site, 1998</p> <p>TH-19 Walsh Test Hole, 1998</p> <p>33 Approximate Property Boundary and ID Number</p> <p>Former Aboveground Storage Tank(s)</p> | <p>Former Underground Storage Tank(s)</p> <p>Underground Storage Tank(s) - Removal Unknown</p> <p>Underground Storage Tank(s)</p> <p>Approximate Location of Former Gas Station</p> <p>Tank Type (UST or AST) Unknown, UST Removal Unknown</p> <p>Former Barrel Storage</p> <p>Walsh Test Hole, 1999</p> | <p>Proposed Monitoring Well</p> <p>Walsh Monitoring Well (1991) Sampling Location</p> <p>Walsh Test Hole, 1998</p> <p>PCE Concentration</p> <p>Aguirre Geotechnical Hole (1994, 1995) Sampling Location</p> | <p>Proposed sidewalks and riprap</p> <p>Proposed curbs and gutters</p> <p>Proposed retaining walls</p> <p>Proposed storm and/or sanitary sewers</p> <p>Proposed gas</p> <p>Proposed telephone</p> <p>Proposed water line</p> |
|---|--|--|---|--|

**Walsh**  
Environmental Scientists and Engineers, Inc.

PCE Concentrations  
In Ground Water

Job 3026-010	Date 9/99	Figure 3
--------------	-----------	----------

3026-010; TH-26 25-27' bgs

**Client Sample ID**

Sample Tag No: 58376

Lab Sample ID: S-99-6-35-2

**Matrix:** Soil

Date Sampled: 06/17/99

Data Filename: VOAA3404.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units:  $\mu\text{g/Kg}$ 

Dilution Factor: 1

[illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

**Analyst:**

Alpha Beta



Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; TH-26 0-1' bgs**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-35-3  
Soil  
Number: 59673  
Sampled: 6/16/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 7/1/99  
Units: mg/Kg

Extractables Dilution Factor: 1

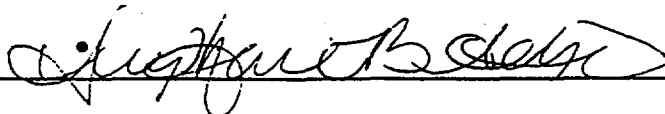
	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA	11	3	

Surrogate Compound	%Recovery
(SS) o-Terphenyl	115 %

## Qualifiers:

- "U" Indicates compound was searched for and not detected at or above the method detection limit.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- " \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_



**Walsh**

Scientists and Engineers, Inc.

3050\_G: Acid Digestion of Sediments, Sludges, and Soils      METHOD: 3050A  
for GFAA Metals

AS\_GTS: ARSENIC, Total (GFAA)      METHOD: 7060

PB\_GTS: LEAD, Total (GFAA)      METHOD: 7421

PMOIST: PERCENT MOISTURE      METHOD: ASTM D2216

# Petroleum Hydrocarbons Report

Page 1 of 1

3026-010; TH-26 25-27' bgs

Client Sample ID

ORIGINAL

EPA Method: Mod. 8100  
Lab Sample ID: 99-6-35-2  
Matrix: Soil  
Tag Number: 58025  
Date Sampled: 6/17/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 7/1/99  
Units: mg/Kg

Extractables Dilution Factor: 1

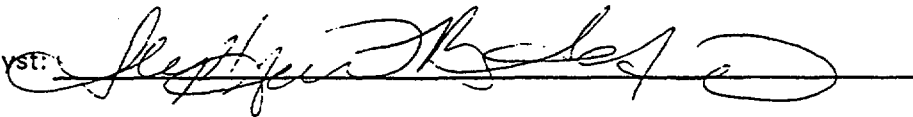
Analyte	CAS Number	Concentration	Detection Limits	Qualifier
Total Extractable Hydrocarbons	NA		3	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	108 %

## Qualifiers:

- "U" Indicates compound was searched for and not detected at or above the method detection limit.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- " \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst:



 **Walsh**

Environmental Scientists and Engineers, Inc.



es were prepared and analyzed according to methods outlined in the  
wing references:

Test Methods for Evaluating Solid Waste, USEPA SW-846, Third Edition,  
Revision 4, December 1996.

Standard Method for Laboratory Determination of Water (Moisture) Content of  
Soil, Rock, and Soil-Aggregate Mixtures, ASTM D 2216-80, July 1980.

ems encountered with the analyses are discussed in the following narrative.

atrix spike recovery for the batched lead quality control sample was below  
ower control limit. All other quality assurance analyses were acceptable.



an Analytica Group company

Walsh Environmental Inc.  
4888 Pearl East Circle #108  
Boulder CO 80301

Attn: Stan Spencer

325 Interlocken Parkway  
Suite 200  
Broomfield, CO 80021  
(303) 469-8868  
(800) 873-8707  
FAX: (303) 469-5254

Order #: 99-06-226  
Date: 06/27/99 16:10  
Work ID: 3026-010 I-70 / BRIGHTON BLVD  
Date Received: 06/18/99  
Date Completed: 06/24/99

#### SAMPLE IDENTIFICATION

<u>Sample</u> <u>Number</u>	<u>Client Description</u>
01	TH-27 TAG #58381

<u>Sample</u> <u>Number</u>	<u>Client Description</u>
--------------------------------	---------------------------

Enclosed are the analytical results for the submitted sample(s). Please review the CASE NARRATIVE for a discussion of any data and/or quality control issues. A listing of data qualifiers and analytical codes is located on the TEST METHODOLOGIES page at the end of the report.

If you have any questions regarding the analyses, please feel free to call.

Sincerely,

Claire K. Toon  
Project Manager

# 99-06-226  
TICA, INC.

Walsh Environmental Inc.  
TEST RESULTS by SAMPLE

Page 3

e: 01A TH-27 TAG #58381

Collected: 06/18/99

Matrix: SOIL

<u>Description</u>	<u>Method</u>	<u>Result</u>	<u>Q</u>	<u>Limit</u>	<u>Units</u>	<u>Analyzed</u>
ic, Total	SW 7060	4.7	DS	0.47	mg/Kg-DRY	06/23/99
Total	SW 7421	68	D	2.4	mg/Kg-DRY	06/23/99
nt Moisture	ASTM D2216	14.9		0.1	WT%	06/23/99

THE FOLLOWING CODES APPLY TO THE ANALYTICAL REPORT

RESULT field...

ND = not detected at the reported limit

NA = analyte not applicable (see case narrative/methods for discussion)

(qualifier) field...

GENERAL:

\* = Recovery or %RPD outside method specifications

H = value is estimated due to analysis run outside EPA holding times

E = reported concentration is above the instrument calibration range

D = analyte was diluted to bring within instrument calibration range or  
to remove matrix interferences

ORGANIC ANALYSIS DATA QUALIFIERS:

B = analyte was detected in the laboratory method blank

J = analyte was detected above the instrument detection limit (IDL)  
but below the analytical reporting limit (CRDL)

INORGANIC ANALYSIS DATA QUALIFIERS:

B = analyte was detected above the instrument detection limit (IDL)  
but below the analytical reporting limit (CRDL)

W = post digestion spike did not meet criteria (85-115%)

S = reported value determined by the Method of Standard Additions

_G:	Acid Digestion of Sediments, Sludges, and Soils for GFAA Metals	METHOD: 3050A
TS:	ARSENIC, Total (GFAA)	METHOD: 7060
TS:	LEAD, Total (GFAA)	METHOD: 7421
ST:	PERCENT MOISTURE	METHOD: ASTM D2216

QA/QC REPORT  
METHOD BLANK SUMMARY  
06/27/99

PAGE: 1  
ORDER#: 9906226

T: JP\_WALSH

ALYTE	UNITS	ANAL DATE	RESULT	LIMIT
SENIC, Total	mg/Kg	06/23/99	ND	0.20
AD, Total	mg/Kg	06/23/99	ND	0.10

METHOD BLANK SPIKE SUMMARY

ALYTE	UNITS	ANAL DATE	RESULT	LIMIT	SPIKE	REF VAL	%REC FLAG	QC SPECS	
								LOW	UPPER
SENIC, Total	mg/Kg	06/23/99	3.8	0.20	4.0	ND	95.0	80	120
AD, Total	mg/Kg	06/23/99	1.9	0.10	2.0	ND	95.0	80	120

MATRIX SPIKE SUMMARY

ALYTE	UNITS	ANAL DATE	RESULT	LIMIT	SPIKE	REF VAL	%REC FLAG	QC SPECS	
								LOW	UPPER
SENIC, Total	mg/Kg	06/23/99	6.9	0.20	4.0	3.8	77.5	70	130
AD, Total	mg/Kg	06/23/99	5.8	0.20	2.0	5.0	40.0	70	130

SAMPLE DUPLICATE SUMMARY

ALYTE	UNITS	ANAL DATE	RESULT	LIMIT	REF VAL	%RPD FLAG	QC SPECS
							UPPER
SENIC, Total	mg/Kg	06/23/99	3.9	0.20	3.8	2.60	.35
AD, Total	mg/Kg	06/23/99	4.9	0.20	5.0	2.02	35

# 99-06-226  
TICA, INC.

Walsh Environmental Inc.  
DATES REPORT

Page 6

e: 01A TH-27 TAG #58381

Matrix: SOIL

	<u>Method</u>	<u>Collected</u>	<u>Received</u>	<u>TCLP date</u>	<u>Extracted</u>	<u>Analyze</u>
Total	SW 7060	06/18/99	06/18/99	NA	06/21/99	06/23/99
Sal	SW 7421	06/18/99	06/18/99	NA	06/21/99	06/23/99
Moisture	ASTM D2216	06/18/99	06/18/99	NA		06/23/99

**ORIGINAL****3026-010; TH-27 0-1'bgs**

Client Sample ID

Sample Tag No: 58389

Lab Sample ID: S-99-6-37-3

Matrix: Soil

Date Extracted: 6/18/99

Data Filename: BNAB1268.D

Date Analyzed: 06/06/99

Analyst: SBS

Dilution Factor: 1

EPA Method: 8270

Units: µg/Kg

Analyte	CAS Number	Concentration	Quantitation Limits	Qualifier
N-Nitrosodimethylamine	62-75-9		660	U
Phenol	108-95-2		330	U
Bis(2-chloroethyl)ether	111-44-4		330	U
2-Chlorophenol	95-57-8		330	U
1,3-Dichlorobenzene	541-73-1		330	U
1,4-Dichlorobenzene	106-46-7		330	U
Benzyl Alcohol	100-51-6		660	U
1,2-Dichlorobenzene	95-50-1		330	U
2-Methylphenol	95-48-7		330	U
Bis(2-chloroisopropyl)ether	108-60-1		330	U
4-Methylphenol	106-44-5		330	U
N-nitrosodi-n-propylamine	621-64-7		330	U
Hexachloroethane	67-72-1		330	U
Nitrobenzene	98-95-3		330	U
Isophorone	78-59-1		330	U
2-Nitrophenol	88-75-5		330	U
2,4-Dimethylphenol	105-67-9		330	U
Bis(2-chloroethoxy)methane	111-91-1		330	U
Benzoic Acid	65-85-0		1,650	U
2,4-Dichlorophenol	120-83-2		330	U
1,2,4-Trichlorobenzene	120-82-1		330	U
Naphthalene	91-20-3		330	U
4-Chloroaniline	106-47-8		660	U
Hexachlorobutadiene	87-68-3		330	U
4-Chloro-3-methylphenol	59-50-7		660	U
2-Methylnaphthalene	91-57-6		330	U
Hexachlorocyclopentadiene	77-47-4		330	U
2,4,6-Trichlorophenol	88-06-2		330	U
2,4,5-Trichlorophenol	95-95-4		330	U
2-Chloronaphthalene	91-58-7		330	U
2-Nitroaniline	88-74-4		1,650	U
Dimethyl Phthalate	131-11-3		330	U
Acenaphthylene	208-96-8		330	U
3-Nitroaniline	99-09-2		1,650	U
Acenaphthene	83-32-9		330	U
2,4-Dinitrophenol	51-28-5		1,650	U
4-Nitrophenol	100-02-7		1,650	U



Environmental Scientists and Engineers, Inc.



**ORIGINAL****3026-010; TH-27 0-1'bgs**

Lab Sample ID: S-99-6-37-3

Client Sample ID:

Dibenzofuran	132-64-9		330	U
2,4-Dinitrotoluene	606-20-2		330	U
2,6-Dinitrotoluene	121-14-2		330	U
Diethylphthalate	84-66-2		330	U
4-Chlorophenyl phenyl ether	7005-72-3		330	U
Fluorene	86-73-7		330	U
4-Nitroaniline	100-01-6		660	U
4,6-Dinitro-2-methylphenol	534-52-1		1,650	U
N-Nitrosodiphenylamine	86-30-6		330	U
4-Bromophenyl-phenylether	101-55-3		330	U
Hexachlorobenzene	118-74-1		330	U
Pentachlorophenol	87-86-5		1,650	U
Phenanthrene	85-01-8		330	U
Anthracene	120-12-7		330	U
Carbazole	86-74-8		660	U
Di-n-butylphthalate	84-74-2		330	U
Fluoranthene	206-44-0		330	U
Pyrene	129-00-0		330	U
Butylbenzylphthalate	85-68-7		330	U
Benzo(a)anthracene	56-55-2		330	U
Chrysene	218-01-9		330	U
3,3'-Dichlorobenzidine	91-94-1		1,650	U
Bis(2-ethylhexyl)phthalate	117-81-7		330	U
Di-n-octylphthalate	117-84-0		330	U
Benzo(b)fluoranthene	205-99-2		330	U
Benzo(k)fluoranthene	207-8-9		330	U
Benzo(a)pyrene	50-32-8		330	U
Indeno(1,2,3-cd)pyrene	193-39-5		660	U
Dibenz(a,h)anthracene	53-70-3		660	U
Benzo(g,h,i)perylene	191-24-2		660	U

Surrogate Compound	%Rec	Limits (%)
(SS) 2-Fluorophenol	85 %	25 --- 121
(SS) Phenol-d5	70 %	24 --- 113
(SS) Nitrobenzene-d5	83 %	23 --- 120
(SS) 2-Fluorobiphenyl	104 %	30 --- 115
(SS) 2,4,6-Tribromophenol	91 %	19 --- 122
(SS) Terphenyl-d14	141 % *	18 --- 137

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"\*\*\*" Indicates surrogates low due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.



Environmental Scientists and Engineers, Inc.

Units:  $\mu\text{g/Kg}$ [illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

Analyst

Clifford B. Allen



Environmental Scientists and Engineers, Inc.



### CHAIN OF CUSTODY RECORD

Nº 8186

LG N# 7706226  
4888 Pearl East Circle, Suite 108  
Boulder Colorado 80301

[illegible]



Environmental Scientists and Engineers, Inc.

CHAIN OF CUSTODY RECORD

№ 8187

4888 Pearl East Circle, Suite 108

Boulder Colorado 80301

# STANDARD TAT

PM: STAN SPENCER 99-6-37

Proj. No.	Project Name
-----------	--------------

3026-010 CDOT I-70/Brighton Blvd

SAMPLERS: (Signature)

*[Handwritten signature]*

[illegible]

Relinquished by: (Sign.)

Date/Time

Received by: (Sign.)

Relinquished by: (Sign.)

Date/Time

Received by: (Sign.)

<i>[Signature]</i>	6/18/99	1318
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Relinquished by: (s) [Signature] / [Signature]

3026-010; TH-27 5-7' bgs

Client Sample ID

ORIGINAL

Sample Tag No: 58384

Lab Sample ID: S-99-6-37-1

Matrix: Soil

Date Sampled: 6/18/99

Data Filename: VOAA3411.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units: µg/Kg

Dilution Factor: 1

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
Perdifluoromethane	75-71-8		10	UJ
Permethane	74-87-3		10	UJ
Chloride	75-01-4		5	UJ
Permethane	74-97-5		10	UJ
Permethane	75-00-3		10	UJ
Perfluoromethane	75-69-4		10	UJ
Trichloroethene	75-35-4		5	UJ
Vinyl Chloride	75-09-2		5	UJ
1,2-Dichloroethene	156-60-5		5	UJ
Trichloroethane	75-34-3		5	UJ
1,2-Dichloroethane	156-59-2		5	UJ
Trichloropropane	594-20-7		5	UJ
Perchloromethane	74-97-5		5	UJ
Perform	67-66-3		5	UJ
Trichloroethane	71-55-6		5	UJ
Trichloropropene	563-58-6		5	UJ
Per Tetrachloride	56-23-5		5	UJ
Trichloroethane	107-06-2		5	UJ
Perene	71-43-2		5	UJ
Permethane	79-01-6		5	UJ
Trichloropropane	78-87-5		5	UJ
Permethane	74-95-3		5	UJ
Perdichloromethane	75-27-4		5	UJ
1,3-Dichloropropene	10061-02-6		5	UJ
Perene	108-88-3		5	UJ
1,3-Dichloropropene	10061-01-5		5	UJ
Trichloroethane	79-00-5		5	UJ
Perbromoethane	106-93-4		5	UJ
Trichloropropane	142-28-9		5	UJ
Trichloroethene	127-18-4		5	UJ
Perchloromethane	124-48-1		5	UJ
Perbenzene	108-90-7		5	UJ
1,2-Tetrachloroethane	630-20-6		5	UJ
Perbenzene	100-41-4		5	UJ
Per-Xylenes			5	UJ
Perene	106-42-3		5	UJ



Environmental Scientists and Engineers, Inc.

3026-010; TH-27 5-7' bgs

Lab Sample ID: S-99-6-37-1

Client Sample ID:

ene	100-42-5		5	UJ
noform	75-25-2		5	UJ
opylbenzene	98-82-8		5	UJ
,2-Tetrachloroethane	79-34-5		5	UJ
nobenzene	108-86-1		5	UJ
-Trichloropropane	96-18-4		5	UJ
pylbenzene	103-65-1		5	UJ
orotoluene	95-49-8		5	UJ
-Trimethylbenzene	108-67-8		5	UJ
/benzene	98-06-6		5	UJ
-Trimethylbenzene	95-63-6		5	UJ
ylbenzene	135-98-8		5	UJ
ichlorobenzene	541-73-1		5	UJ
propyltoluene	99-87-6		5	UJ
ichlorobenzene	106-46-7		5	UJ
ylbenzene	104-51-8		5	UJ
ichlorobenzene	95-50-1		5	UJ
ibromo-3-chloropropane	96-12-8		5	UJ
Trichlorobenzene	120-82-1		5	UJ
chlorobutadiene	87-68-3		5	UJ
halene	91-20-3		5	UJ
Trichlorobenzene	87-61-6		5	UJ

gate Compound	%Rec	Limits (%)
Dibromofluoromethane	102 %	80 --- 120
toluene-d8	105 %	81 --- 117
p-Bromofluorobenzene	99 %	74 --- 121

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"\*" Indicates surrogates low due to matrix effect.



ental Scientists and Engineers, Inc.

3026-010; TH-27 5-7' bgs

Client Sample ID

Sample Tag No: 58384

Lab Sample ID: S-99-6-37-1

Matrix: Soil

Date Sampled: 06/18/99

Data Filename: VOAA3411.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units:  $\mu\text{g/Kg}$ 

Dilution Factor: 1

[illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

Analyst: Christopher Baker

# Walsh

Scientists and Engineers, Inc.

026-010; TH-27 25-27' bgs

Client Sample ID

ORIGINAL

Sample Tag No: 58385

Lab Sample ID: S-99-6-37-2

Matrix: Soil

Date Sampled: 6/18/99

Data Filename: VOAA3410.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units: µg/Kg

Dilution Factor: 1

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
Acrylonitrile	75-71-8		10	UJ
Chloroethane	74-87-3		10	UJ
Chloride	75-01-4		5	UJ
Chloroethane	74-97-5		10	UJ
Chloroethane	75-00-3		10	UJ
Chlorofluoromethane	75-69-4		10	UJ
Dichloroethene	75-35-4		5	UJ
Dichloroethene Chloride	75-09-2		5	UJ
1,2-Dichloroethene	156-60-5		5	UJ
Dichloroethane	75-34-3		5	UJ
1,2-Dichloroethane	156-59-2		5	UJ
Dichloropropane	594-20-7		5	UJ
Dichloromethane	74-97-5		5	UJ
Form	67-66-3		5	UJ
Trichloroethane	71-55-6		5	UJ
Dichloropropene	563-58-6		5	UJ
Carbon Tetrachloride	56-23-5		5	UJ
Dichloroethane	107-06-2		UJ	UJ
ene	71-43-2		5	UJ
Chloroethene	79-01-6		5	UJ
Dichloropropane	78-87-5		5	UJ
Chloromethane	74-95-3		5	UJ
Dichloromethane	75-27-4		5	UJ
1,3-Dichloropropene	10061-02-6		5	UJ
ene	108-88-3		5	UJ
1,3-Dichloropropene	10061-01-5		5	U
Trichloroethane	79-00-5		5	U
Bromoethane	106-93-4		5	U
Chloropropane	142-28-9		5	U
Chloroethene	127-18-4		5	U
Dichloromethane	124-48-1		5	U
Benzene	108-90-7		5	U
1,2-Tetrachloroethane	630-20-6		5	U
Benzene	100-41-4		5	U
m-Xylenes			5	U
ene	106-42-3		5	U

**Walsh**

Environmental Scientists and Engineers, Inc.



3026-010; TH-27 25-27' bgs

Lab Sample ID: S-99-6-37-2

Client Sample ID:

ene	100-42-5		5	U
noform	75-25-2		5	UJ
propylbenzene	98-82-8		5	UJ
1,2-Tetrachloroethane	79-34-5		5	UJ
nobenzene	108-86-1		5	UJ
-Trichloropropane	96-18-4		5	UJ
propylbenzene	103-65-1		5	UJ
lorotoluene	95-49-8		5	UJ
-Trimethylbenzene	108-67-8		5	UJ
ylbenzene	98-06-6		5	UJ
-Trimethylbenzene	95-63-6		5	UJ
ylbenzene	135-98-8		5	UJ
ichlorobenzene	541-73-1		5	UJ
propyltoluene	99-87-6		5	UJ
ichlorobenzene	106-46-7		5	UJ
ylbenzene	104-51-8		5	UJ
ichlorobenzene	95-50-1		5	UJ
ibromo-3-chloropropane	96-12-8		5	UJ
-Trichlorobenzene	120-82-1		5	UJ
chlorobutadiene	87-68-3		5	UJ
thalene	91-20-3		5	UJ
-Trichlorobenzene	87-61-6		5	UJ

ogate Compound	%Rec	Limits (%)
Dibromofluoromethane	104 %	80 — 120
Toluene-d8	107 %	81 — 117
o-Bromofluorobenzene	105 %	74 — 121

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

\*\*\* Indicates surrogates low due to matrix effect.



mental Scientists and Engineers, Inc.

Client Sample ID  
Sample Tag No: 58387  
Lab Sample ID: S-99-6-37-2  
Matrix: Soil  
Date Sampled: 06/18/99  
Data Filename: VOAA3410.D

EPA Method: 8260  
Date Analyzed: 06/22/99  
Analyst: SBS  
Units: µg/Kg  
Dilution Factor: 1

[illegible]

"T" Indicates compound was tentatively identified by its mass spectrum.  
All tentatively identified compounds are estimated values.

**Analyst:**

Christopher B. Lehn



# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; TH-27 5-7' bgs**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-37-1  
X: Soil  
Number: 58386  
Sampled: 6/18/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 7/1/99  
Units: mg/Kg

Extractables Dilution Factor: 1

te	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		3	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	115 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*[Signature]*

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; TH-27 25-27' bgs**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-37-2  
Soil  
Number: 58387  
Sampled: 6/18/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 7/1/99  
Units: mg/Kg

Extractables Dilution Factor: 1

	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		3	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	69 %

## Qualifiers:

- "U" Indicates compound was searched for and not detected at or above the method detection limit.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- " \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Stephen B. Davis*

**Walsh**

Scientists and Engineers, Inc.

ORIGINAL

026-010; TH-28 1-3' bgs

Client Sample ID

Sample Tag No: 59677

Sample ID: S-99-6-35-6

Matrix: Soil

Date Extracted: 6/25/99

Data Filename: BNAB1270.D

Date Analyzed: 07/06/99

Analyst: SBS

Dilution Factor: 1

EPA Method: 8270

Units: µg/Kg

	CAS Number	Concentration	Quantitation Limits	Qualifier
dimethylamine	62-75-9		660	U
	108-95-2		330	U
diethylether	111-44-4		330	U
phenol	95-57-8		330	U
chlorobenzene	541-73-1		330	U
chlorobenzene	106-46-7		330	U
alcohol	100-51-6		660	U
chlorobenzene	95-50-1		330	U
phenol	95-48-7		330	U
diisopropylether	108-60-1		330	U
phenol	106-44-5		330	U
di-n-propylamine	621-64-7		330	U
proethane	67-72-1		330	U
zene	98-95-3		330	U
ne	78-59-1		330	U
enol	88-75-5		330	U
thylphenol	105-67-9		330	U
proethoxy)methane	111-91-1		330	U
Acid	65-85-0		1,650	U
rophenol	120-83-2		330	U
chlorobenzene	120-82-1		330	U
ene	91-20-3		330	U
aniline	106-47-8		660	U
robutadiene	87-68-3		330	U
3-methylphenol	59-50-7		660	U
naphthalene	91-57-6		330	U
rocyclopentadiene	77-47-4		330	U
chlorophenol	88-06-2		330	U
chlorophenol	95-95-4		330	U
naphthalene	91-58-7		330	U
line	88-74-4		1,650	U
Phthalate	131-11-3		330	U
nylene	208-96-8		330	U
line	99-09-2		1,650	U
hene	83-32-9		330	U
ophenol	51-28-5		1,650	U
enol	100-02-7		1,650	U

Walsh

Scientists and Engineers, Inc.

026-010; TH-28 1-3' bgs

Lab Sample ID: S-99-6-35-6

Client Sample ID:

zofuran	132-64-9		330	U
nitrotoluene	606-20-2		330	U
nitrotoluene	121-14-2		330	U
lphthalate	84-66-2		330	U
rophenyl phenyl ether	7005-72-3		330	U
ne	86-73-7		330	U
aniline	100-01-6		660	U
nitro-2-methylphenol	534-52-1		1,650	U
osodiphenylamine	86-30-6		330	U
nophenyl-phenylether	101-55-3		330	U
olorobenzene	118-74-1		330	U
hlorophenol	87-86-5		1,650	U
anthrene	85-01-8	340	330	
cene	120-12-7		330	U
ole	86-74-8		660	U
tylphthalate	84-74-2		330	U
nthene	206-44-0	560	330	
	129-00-0	920	330	
enzylphthalate	85-68-7		330	U
a)anthracene	56-55-2		330	U
ne	218-01-9	370	330	
chlorobenzidine	91-94-1		1,650	U
thylhexyl)phthalate	117-81-7		330	U
tylphthalate	117-84-0		330	UJ
b)fluoranthene	205-99-2	430	330	J
k)fluoranthene	207-8-9		330	UJ
a)pyrene	50-32-8		330	UJ
1,2,3-cd)pyrene	193-39-5		660	UJ
(a,h)anthracene	53-70-3		660	UJ
g,h,i)perylene	191-24-2		660	UJ

ate Compound	%Rec	Limits (%)
Fluorophenol	58 %	25 --- 121
enol-d5	48 %	24 --- 113
robenzene-d5	70 %	23 --- 120
Fluorobiphenyl	87 %	30 --- 115
1,6-Tribromophenol	83 %	19 --- 122
rphenyl-d14	129 %	18 --- 137

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"" Indicates surrogates low due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.

**Walsh**

Anal Scientists and Engineers, Inc.

ORIGINAL

3026-010; TH-28 1-3' bgs

Client Sample ID

Sample Tag No: 59677

Lab Sample ID: S-99-6-35-6

Matrix: Soil

Data Filename: BNAB1270.D

Date Analyzed: 07/06/99

Analyst: SBS

Dilution Factor: 1

Method: 8270

Units: µg/Kg

tatively Identified Compound	Concentration	Qualifier
nown Organic Compound	140	TJB
nown Organic Compound	190	TJB
nown Organic Compound	1800	TJB
yclohexen-1-one	170	TJ
nown Substituted Alkane	410	TJ
nown Organic Compound	220	TJ
nown Organic Compound	480	TJ
nown Substituted Alkene	170	TJ
ecular Sulfur	1900	TJ
nown Organic Compound	770	TJ

Qualifier:

"T" Indicates compound was tentatively identified by its mass spectrum.


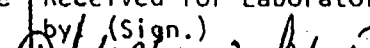
All tentatively identified compounds are estimated values.

Analyst

*Stephen B. DeGru*

Environmental Scientists and Engineers, Inc.

*[Handwritten signature]*

Relinquished by: (Sign.) 	Date/Time 6/17/99 1624	Received by: (Sign.)	Relinquished by: (Sign.)	Date/Time	Received by: (Sign.)
Relinquished by: (Sign.)	Date/Time	Received by: (Sign.)	Relinquished by: (Sign.)	Date/Time	Received by: (Sign.)
Relinquished by: (Sign.)	Date/Time	Received for Laboratory by (Sign.) 	Date/Time 6/17/99 1624	Remarks:	

Distribution: Original accompanies element 1208011 (1/1/1997)



**3026-010; TH-28 10-12' bgs**

Client Sample ID

**ORIGINAL**

Sample Tag No: 58027

Lab Sample ID: S-99-6-35-4

Matrix: Soil

Date Sampled: 6/17/99

Data Filename: VOAA3405.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units: µg/Kg

Dilution Factor: 1

Analyte	CAS Number	Concentration	Quantitation Limits	Qualifier
Dichlorodifluoromethane	75-71-8		10	UJ
Chloromethane	74-87-3		10	UJ
Vinyl Chloride	75-01-4		5	UJ
Bromomethane	74-97-5		10	UJ
Chloroethane	75-00-3		10	UJ
Trichlorofluoromethane	75-69-4		10	UJ
1,1-Dichloroethene	75-35-4		5	UJ
Methylene Chloride	75-09-2		5	UJ
trans-1,2-Dichloroethene	156-60-5		5	UJ
1,1-Dichloroethane	75-34-3		5	UJ
cis-1,2-Dichloroethene	156-59-2		5	UJ
2,2-Dichloropropane	594-20-7		5	UJ
Bromochloromethane	74-97-5		5	UJ
Chloroform	67-66-3		5	UJ
1,1,1-Trichloroethane	71-55-6		5	UJ
1,1-Dichloropropene	563-58-6		5	UJ
Carbon Tetrachloride	56-23-5		5	UJ
1,2-Dichloroethane	107-06-2		5	UJ
Benzene	71-43-2		5	UJ
Trichloroethene	79-01-6		5	UJ
1,2-Dichloropropane	78-87-5		5	UJ
Dibromomethane	74-95-3		5	UJ
Bromodichloromethane	75-27-4		5	UJ
cis-1,3-Dichloropropene	10061-02-6		5	UJ
Toluene	108-88-3		5	UJ
trans-1,3-Dichloropropene	10061-01-5		5	UJ
1,1,2-Trichloroethane	79-00-5		5	UJ
1,2-Dibromoethane	106-93-4		5	UJ
1,3-Dichloropropane	142-28-9		5	UJ
Tetrachloroethene	127-18-4		5	UJ
Dibromochloromethane	124-48-1		5	UJ
Chlorobenzene	108-90-7		5	UJ
1,1,1,2-Tetrachloroethane	630-20-6		5	UJ
Ethylbenzene	100-41-4		5	UJ
m & p-Xylenes			5	UJ
o-Xylene	106-42-3		5	UJ



Environmental Scientists and Engineers, Inc.

026-010; TH-28 25-27' bgs

Client Sample ID

ORIGINAL

Sample Tag No: 58026

Lab Sample ID: S-99-6-35-5

Matrix: Soil

Date Sampled: 6/17/99

Data Filename: VOAA3436.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units: µg/Kg

Dilution Factor: 1

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
Fluoromethane	75-71-8		10	U
Methane	74-87-3		10	U
Chloride	75-01-4		5	U
Methane	74-97-5		10	U
Ethane	75-00-3		10	U
Fluoromethane	75-69-4		10	U
Chloroethene	75-35-4		5	U
Ene Chloride	75-09-2		5	U
1,2-Dichloroethene	156-60-5		5	U
Chloroethane	75-34-3		5	U
1,1-Dichloroethene	156-59-2		5	U
Chloropropane	594-20-7		5	U
Chloromethane	74-97-5		5	U
Form	67-66-3		5	U
Trichloroethane	71-55-6		5	U
Chloropropene	563-58-6		5	U
1,1,1-Tetrachloride	56-23-5		5	U
Chloroethane	107-06-2		5	U
Ene	71-43-2		5	U
Chloroethene	79-01-6		5	U
Chloropropane	78-87-5		5	U
Methane	74-95-3		5	U
Dichloromethane	75-27-4		5	U
Dichloropropene	10061-02-6		5	U
Ene	108-88-3		5	U
1,3-Dichloropropene	10061-01-5		5	U
Trichloroethane	79-00-5		5	U
Chloroethane	106-93-4		5	U
Chloropropane	142-28-9		5	U
Chloroethene	127-18-4		5	U
Chloromethane	124-48-1		5	U
Benzene	108-90-7		5	U
Tetrachloroethane	630-20-6		5	U
Benzene	100-41-4		5	U
Ylenes			5	U
Ene	106-42-3		5	U

**Walsh**

Anal Scientists and Engineers, Inc.

026-010; TH-28 25-27' bgs

**Client Sample ID**

Sample Tag No: 58026

Lab Sample ID: S-99-6-35-5

Matrix: Soil

Date Sampled: 06/17/99

Data Filename: VOAA3406.D

EPA Method: 8260

Date Analyzed: 06/22/99

Analyst: SBS

Units:  $\mu\text{g/Kg}$ 

Dilution Factor: 1

[illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

Analyst: Jeffrey D. Bole



**ital Scientists and Engineers, Inc.**

3026-010; TH-28 25-27' bgs

Lab Sample ID: S-99-6-35-5

Client Sample ID:

ene	100-42-5		5	U
noform	75-25-2		5	U
ropylbenzene	98-82-8		5	U
2,2-Tetrachloroethane	79-34-5		5	U
nobenzene	108-86-1		5	U
3-Trichloropropane	96-18-4		5	U
opylbenzene	103-65-1		5	U
lorotoluene	95-49-8		5	U
-Trimethylbenzene	108-67-8		5	U
ylbenzene	98-06-6		5	U
-Trimethylbenzene	95-63-6		5	U
ylbenzene	135-98-8		5	U
ichlorobenzene	541-73-1		5	U
propyltoluene	99-87-6		5	U
ichlorobenzene	106-46-7		5	U
ylbenzene	104-51-8		5	U
ichlorobenzene	95-50-1		5	U
ibromo-3-chloropropane	96-12-8		5	U
Trichlorobenzene	120-82-1		5	U
chlorobutadiene	87-68-3		5	U
halene	91-20-3		5	U
Trichlorobenzene	87-61-6		5	U

gate Compound	%Rec	Limits (%)
Dibromofluoromethane	99 %	80 --- 120
toluene-d8	103 %	81 --- 117
p-Bromofluorobenzene	103 %	74 --- 121

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"\*\*" Indicates surrogates low due to matrix effect.



Walsh Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; TH-28 10-12' bgs**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-35-4  
Soil  
Number: 58383  
Sampled: 6/17/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 7/1/99  
Units: mg/Kg

Extractables Dilution Factor: 1

	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA	6.0	3	

Surrogate Compound	%Recovery
(SS) o-Terphenyl	102 %

## Qualifiers:

- "U" Indicates compound was searched for and not detected at or above the method detection limit.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- " \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Christopher D. Bolognino*

**Walsh**  
Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; TH-28 25-27' bgs**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-35-5  
Soil  
Number: 58028  
Sampled: 6/17/99

Analyst: SBS  
Date Extracted: 6/29/99  
Extractables Date Analyzed: 6/30/99  
Units: mg/Kg

Extractables Dilution Factor: 1

	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		3	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	80 %

## Qualifiers:

- "U" Indicates compound was searched for and not detected at or above the method detection limit.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- " \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*[Signature]*

**Walsh**

Environmental Scientists and Engineers, Inc.

# Volatile Organic Compounds Report

Page 1 of 3

3026-010; TH-7

## ORIGINAL

Client Sample ID

Sample Tag No.: 59012, 59013

Lab Sample ID: W-99-3-30-4

Matrix: Water

Data Filename: VOAA3213.D

EPA Method: 8260

Date Sampled: 03/23/99

Date Analyzed: 03/24/99

Analyst: SBS

Units:  $\mu\text{g/L}$

Dilution Factor: 1

Analyte	CAS Number	Concentration	Quantitation Limits	Qualifier
Dichlorodifluoromethane	75-71-8		10	U
Chloromethane	74-87-3		10	U
Vinyl Chloride	75-01-4		5	U
Bromomethane	74-97-5		10	U
Chloroethane	75-00-3		10	U
Trichlorofluoromethane	75-69-4		10	U
1,1-Dichloroethene	75-35-4		5	U
Methylene Chloride	75-09-2		5	U
trans-1,2-Dichloroethene	156-60-5		5	U
1,1-Dichloroethane	75-34-3		5	U
cis-1,2-Dichloroethene	156-59-2		5	U
2,2-Dichloropropane	594-20-7		5	U
Bromochloromethane	74-97-5		5	U
Chloroform	67-66-3		5	U
1,1,1-Trichloroethane	71-55-6		5	U
1,1-Dichloropropene	563-58-6		5	U
Carbon Tetrachloride	56-23-5		5	U
1,2-Dichloroethane	107-06-2		5	U
Benzene	71-43-2		5	U
Trichloroethene	79-01-6		5	U
1,2-Dichloropropane	78-87-5		5	U
Dibromomethane	74-95-3		5	U
Bromodichloromethane	75-27-4		5	U
trans-1,3-Dichloropropene	10061-02-6		5	U
Toluene	108-88-3		5	U
cis-1,3-Dichloropropene	10061-01-5		5	U
1,1,2-Trichloroethane	79-00-5		5	U
1,2-Dibromoethane	106-93-4		5	U
1,3-Dichloropropane	142-28-9		5	U
Tetrachloroethene	127-18-4	61	25	
Dibromochloromethane	124-48-1		5	U
Chlorobenzene	108-90-7		5	U
1,1,1,2-Tetrachloroethane	630-20-6		5	U
Ethylbenzene	100-41-4		5	U
m & p-Xylenes			5	U



Environmental Scientists and Engineers, Inc.

# Volatile Organic Compounds Report

Page 2 of 3

3026-010; TH-7

Lab Sample ID: W-99-3-30-4

## Client Sample ID

o-Xylene	106-42-3	5	U
Styrene	100-42-5	5	U
Bromoform	75-25-2	5	U
Isopropylbenzene	98-82-8	5	U
1,1,2,2-Tetrachloroethane	79-34-5	5	U
Bromobenzene	108-86-1	5	U
1,2,3-Trichloropropane	96-18-4	5	U
n-Propylbenzene	103-65-1	5	U
2-Chlorotoluene	95-49-8	5	U
1,3,5-Trimethylbenzene	108-67-8	5	U
4-Chlorotoluene	106-43-4	5	U
t-Butylbenzene	98-06-6	5	U
1,2,4-Trimethylbenzene	95-63-6	5	U
s-Butylbenzene	135-98-8	5	U
1,3-Dichlorobenzene	541-73-1	5	U
p-Isopropyltoluene	99-87-6	5	U
1,4-Dichlorobenzene	106-46-7	5	U
n-Butylbenzene	104-51-8	5	U
1,2-Dichlorobenzene	95-50-1	5	U
1,2-Dibromo-3-chloropropane	96-12-8	5	U
1,2,4-Trichlorobenzene	120-82-1	5	U
Hexachlorobutadiene	87-68-3	5	U
Naphthalene	91-20-3	10	U
1,2,3-Trichlorobenzene	87-61-6	5	U

Surrogate Compound	%Rec	Recovery Limits (%)		
(SS) Dibromofluoromethane	98 %	86	---	118
(SS) Toluene-d8	101 %	88	---	110
(SS) p-Bromofluorobenzene	100 %	86	---	116

### Qualifiers:

- "U" Indicates compound was searched for and not detected.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- "D" Indicates compound was run at a dilution.
- "\*" Indicates surrogate recovery is not within method limits due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.



## Page 3 of 3

Client Sample ID

Data Filename: VOAA3213.D

Units:  $\mu\text{g/L}$ 

Tentatively Identified Compound	Concentration	Qualifier
No Tentatively Identified Compounds were detected.		

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

**Analyst:**

Stephen B. Kiser

# Petroleum Hydrocarbons Report

Page 1 of 1

3026-010; TH-7

ORIGINAL

Method: mod. 8015/mod. 8100  
Sample ID: 99-3-30-4  
Matrix: Water  
Sample Number: 59009, 10, 11  
Date Sampled: 03/23/99

Analyst: DPD  
Volatiles Date Analyzed: 03/24/99  
Date Extracted: 03/24/99  
Extractables Date Analyzed: 03/25/99  
Units:  $\mu\text{g/L}$

Volatiles Dilution Factor: 1

Extractables Dilution Factor: 1

Analyte	CAS Number	Concentration	Reporting Limits	Qualifier
Total Volatile Hydrocarbons	NA		500	U
Total Extractable Hydrocarbons	NA	1300	1000	

Surrogate Compound	%Recovery
(SS) a,a,a-Trifluorotoluene	75 %
(SS) Fluorobenzene	115 %
(SS) o-Terphenyl	56 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank and has been corrected.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Dangming Dai*

**Walsh**

Environmental Scientists and Engineers, Inc.

# Volatile Organic Compounds Report

Page 1 of 3

3026-010; TH-19

ORIGINAL

Client Sample ID

Sample Tag No.: 59007, 59008

Lab Sample ID: W-99-3-30-3

Matrix: Water

Data Filename: VOAA3212.D

EPA Method: 8260

Date Sampled: 03/23/99

Date Analyzed: 03/24/99

Analyst: SBS

Units: µg/L

Dilution Factor: 1

anlyte	CAS Number	Concentration	Quantitation Limits	Qualifier
chlorodifluoromethane	75-71-8		10	U
loromethane	74-87-3		10	U
nyl Chloride	75-01-4		5	U
omomethane	74-97-5		10	U
loroethane	75-00-3		10	U
chlorofluoromethane	75-69-4		10	U
1-Dichloroethene	75-35-4		5	U
ethylene Chloride	75-09-2		5	U
ns-1,2-Dichloroethene	156-60-5		5	U
1-Dichloroethane	75-34-3		5	U
-1,2-Dichloroethene	156-59-2		5	U
2-Dichloropropane	594-20-7		5	U
omochloromethane	74-97-5		5	U
loroform	67-66-3		5	U
,1-Trichloroethane	71-55-6		5	U
-Dichloropropene	563-58-6		5	U
arbon Tetrachloride	56-23-5		5	U
2-Dichloroethane	107-06-2		5	U
azene	71-43-2		5	U
chloroethene	79-01-6		5	U
-Dichloropropane	78-87-5		5	U
romomethane	74-95-3		5	U
modichloromethane	75-27-4		5	U
ns-1,3-Dichloropropene	10061-02-6		5	U
uene	108-88-3		5	U
1,3-Dichloropropene	10061-01-5		5	U
,2-Trichloroethane	79-00-5		5	U
-Dibromoethane	106-93-4		5	U
-Dichloropropane	142-28-9		5	U
rachloroethene	127-18-4	160	25	
romochloromethane	124-48-1		5	U
robenzene	108-90-7		5	U
,1,2-Tetrachloroethane	630-20-6		5	U
ylbenzene	100-41-4		5	U
p-Xylenes			5	U

# Volatile Organic Compounds Report

Page 2 of 3

3026-010; TH-19

Lab Sample ID: W-99-3-30-3

## Client Sample ID

o-Xylene	106-42-3	5	U
Styrene	100-42-5	5	U
Bromoform	75-25-2	5	U
Isopropylbenzene	98-82-8	5	U
1,1,2,2-Tetrachloroethane	79-34-5	5	U
Bromobenzene	108-86-1	5	U
1,2,3-Trichloropropane	96-18-4	5	U
n-Propylbenzene	103-65-1	5	U
2-Chlorotoluene	95-49-8	5	U
1,3,5-Trimethylbenzene	108-67-8	5	U
4-Chlorotoluene	106-43-4	5	U
t-Butylbenzene	98-06-6	5	U
1,2,4-Trimethylbenzene	95-63-6	5	U
s-Butylbenzene	135-98-8	5	U
1,3-Dichlorobenzene	541-73-1	5	U
p-Isopropyltoluene	99-87-6	5	U
1,4-Dichlorobenzene	106-46-7	5	U
n-Butylbenzene	104-51-8	5	U
1,2-Dichlorobenzene	95-50-1	5	U
1,2-Dibromo-3-chloropropane	96-12-8	5	U
1,2,4-Trichlorobenzene	120-82-1	5	U
Hexachlorobutadiene	87-68-3	5	U
Naphthalene	91-20-3	10	U
1,2,3-Trichlorobenzene	87-61-6	5	U

Surrogate Compound	%Rec	Recovery Limits (%)		
(SS) Dibromofluoromethane	100 %	86	---	118
(SS) Toluene-d8	101 %	88	---	110
(SS) p-Bromofluorobenzene	101 %	86	---	116

### Qualifiers:

"U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"D" Indicates compound was run at a dilution.

"\*\*" Indicates surrogate recovery is not within method limits due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.

## Page 3 of 3

**Client Sample ID**

Data Filename: VOAA3212.D

Units:  $\mu\text{g/L}$ 

Qualifier:

"T" Indicates compound was tentatively identified by its mass spectrum.  
All tentatively identified compounds are estimated values.

**Analyst:**



**Walsh**  
Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

3026-010; TH-19

ORIGINAL

EPA Method: mod. 8015/mod. 8100  
Lab Sample ID: 99-3-30-3  
Matrix: Water  
Tag Number: 59004, 05, 06  
Date Sampled: 03/23/99

Analyst: DPD  
Volatiles Date Analyzed: 03/24/99  
Date Extracted: 03/24/99  
Extractables Date Analyzed: 03/25/99  
Units:  $\mu\text{g/L}$

Volatiles Dilution Factor: 1

Extractables Dilution Factor: 1

Analyte	CAS Number	Concentration	Reporting Limits	Qualifier
Total Volatile Hydrocarbons	NA		500	U
Total Extractable Hydrocarbons	NA		1000	U

Surrogate Compound	% Recovery
(SS) a,a,a-Trifluorotoluene	71 %
(SS) Fluorobenzene	111 %
(SS) o-Terphenyl	64 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank and has been corrected.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Dongming Dai*

# Volatile Organic Compounds Report

Page 1 of 3

3026-010; TH-24

## ORIGINAL

Client Sample ID

Sample Tag No.: 58997, 58998

Lab Sample ID: W-99-3-30-1

Matrix: Water

Data Filename: VOAA3214.D

VOAA3245.D

EPA Method: 8260

Date Sampled: 03/23/99

Date Analyzed: 3/24/99

Analyst: SBS

Units:  $\mu\text{g/L}$

Dilution Factor: 5, 2

Analyte	CAS Number	Concentration	Quantitation Limits	Qualifier
Dichlorodifluoromethane	75-71-8		20	U
Chloromethane	74-87-3		20	U
Vinyl Chloride	75-01-4		10	U
Bromomethane	74-97-5		20	U
Chloroethane	75-00-3		20	U
Trichlorofluoromethane	75-69-4		20	U
1,1-Dichloroethene	75-35-4		10	U
Methylene Chloride	75-09-2		10	U
trans-1,2-Dichloroethene	156-60-5		10	U
1,1-Dichloroethane	75-34-3		10	U
cis-1,2-Dichloroethene	156-59-2		10	U
2,2-Dichloropropane	594-20-7		10	U
Bromochloromethane	74-97-5		10	U
Chloroform	67-66-3		10	U
1,1,1-Trichloroethane	71-55-6		10	U
1,1-Dichloropropene	563-58-6		10	U
Carbon Tetrachloride	56-23-5		10	U
1,2-Dichloroethane	107-06-2		10	U
Benzene	71-43-2		10	U
Trichloroethene	79-01-6		10	U
1,2-Dichloropropane	78-87-5		10	U
Dibromomethane	74-95-3		10	U
Bromodichloromethane	75-27-4		10	U
trans-1,3-Dichloropropene	10061-02-6		10	U
Toluene	108-88-3		10	U
cis-1,3-Dichloropropene	10061-01-5		10	U
1,1,2-Trichloroethane	79-00-5		10	U
1,2-Dibromoethane	106-93-4		10	U
1,3-Dichloropropane	142-28-9		10	U
Tetrachloroethene	127-18-4	480	25	D
Dibromochloromethane	124-48-1		10	U
Chlorobenzene	108-90-7		10	U
1,1,1,2-Tetrachloroethane	630-20-6		10	U
Ethylbenzene	100-41-4		10	U
m & p-Xylenes			10	U

# Volatile Organic Compounds Report

Page 2 of 3

3026-010; TH-24

Lab Sample ID: W-99-3-30-1

## Client Sample ID

o-Xylene	106-42-3	10	U
Styrene	100-42-5	10	U
Bromoform	75-25-2	10	U
Isopropylbenzene	98-82-8	10	U
1,1,2,2-Tetrachloroethane	79-34-5	10	U
Bromobenzene	108-86-1	10	U
1,2,3-Trichloropropane	96-18-4	10	U
n-Propylbenzene	103-65-1	10	U
2-Chlorotoluene	95-49-8	10	U
1,3,5-Trimethylbenzene	108-67-8	10	U
4-Chlorotoluene	106-43-4	10	U
t-Butylbenzene	98-06-6	10	U
1,2,4-Trimethylbenzene	95-63-6	10	U
s-Butylbenzene	135-98-8	10	U
1,3-Dichlorobenzene	541-73-1	10	U
p-Isopropyltoluene	99-87-6	10	U
1,4-Dichlorobenzene	106-46-7	10	U
n-Butylbenzene	104-51-8	10	U
1,2-Dichlorobenzene	95-50-1	10	U
1,2-Dibromo-3-chloropropane	96-12-8	10	U
1,2,4-Trichlorobenzene	120-82-1	10	U
Hexachlorobutadiene	87-68-3	10	U
Naphthalene	91-20-3	20	U
1,2,3-Trichlorobenzene	87-61-6	10	U

Surrogate Compound	%Rec	Recovery Limits (%)
(SS) Dibromofluoromethane	101 %	86 --- 118
(SS) Toluene-d8	101 %	88 --- 110
(SS) p-Bromofluorobenzene	104 %	86 --- 116

### Qualifiers:

- "U" Indicates compound was searched for and not detected.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- "D" Indicates compound was run at a dilution.
- "\*" Indicates surrogate recovery is not within method limits due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.



## Page 3 of 3

**Client Sample ID**

Units:  $\mu\text{g/L}$ 

**Qualifier:**

All tentatively identified compounds are estimated values.

**Analyst:**

Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**ORIGINAL**

**3026-010; TH-24**

PA Method: mod. 8015/mod. 8100  
 Sample ID: 99-3-30-1  
 Matrix: Water  
 Log Number: 58994, 95, 96  
 Date Sampled: 03/23/99

Analyst: DPD  
 Volatiles Date Analyzed: 03/24/99  
 Date Extracted: 03/24/99  
 Extractables Date Analyzed: 03/25/99  
 Units: µg/L

Volatiles Dilution Factor: 1

Extractables Dilution Factor: 1

Analyste	CAS Number	Concentration	Reporting Limits	Qualifier
Total Volatile Hydrocarbons	NA	600	500	
Total Extractable Hydrocarbons	NA		1000	U

Surrogate Compound	%Recovery
(SS) a,a,a-Trifluorotoluene	77 %
(SS) Fluorobenzene	125 %
(SS) o-Terphenyl	39 % *

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank and has been corrected.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst:

*Dongming Dai*

# Volatile Organic Compounds Report

Page 1 of 3

3026-010; TH-25

Client Sample ID

ORIGINAL

Sample Tag No.: 59002, 59003

Lab Sample ID: W-99-3-30-2

Matrix: Water

Data Filename: VOAA3238.D

EPA Method: 8260

Date Sampled: 03/23/99

Date Analyzed: 03/26/99

Analyst: SBS

Units: µg/L

Dilution Factor: 1

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
Perfluoromethane	75-71-8		10	U
Methane	74-87-3		10	U
Chloride	75-01-4		5	U
Methane	74-97-5		10	U
Ethane	75-00-3		10	U
Perfluoromethane	75-69-4		10	U
Chloroethene	75-35-4		5	U
Ene Chloride	75-09-2		5	U
1,2-Dichloroethene	156-60-5		5	U
Chloroethane	75-34-3		5	U
1,1-Dichloroethene	156-59-2		5	U
Chloropropane	594-20-7		5	U
Chloromethane	74-97-5		5	U
Form	67-66-3		5	U
Trichloroethane	71-55-6	6.9	5	
Chloropropene	563-58-6		5	U
Tetrachloride	56-23-5		5	U
Chloroethane	107-06-2		5	U
	71-43-2		5	U
Chloroethene	79-01-6		5	U
Chloropropane	78-87-5		5	U
Chloromethane	74-95-3		5	U
1,1-Dichloromethane	75-27-4		5	U
1,3-Dichloropropene	10061-02-6		5	U
	108-88-3		5	U
Dichloropropene	10061-01-5		5	U
1,1-Dichloroethane	79-00-5		5	U
Chloroethane	106-93-4		5	U
Chloropropane	142-28-9		5	U
Chloroethene	127-18-4	23	5	
Chloromethane	124-48-1		5	U
Benzene	108-90-7		5	U
Tetrachloroethane	630-20-6		5	U
Benzene	100-41-4		5	U
Arenes			5	U

# Volatile Organic Compounds Report

Page 2 of 3

3026-010; TH-25

Lab Sample ID: W-99-3-30-2

## Client Sample ID

ene	106-42-3	5	U
ane	100-42-5	5	U
oform	75-25-2	5	U
opylbenzene	98-82-8	5	U
2,2-Tetrachloroethane	79-34-5	5	U
obenzene	108-86-1	5	U
3-Trichloropropane	96-18-4	5	U
pylbenzene	103-65-1	5	U
lorotoluene	95-49-8	5	U
-Trimethylbenzene	108-67-8	5	U
orotoluene	106-43-4	5	U
ylbenzene	98-06-6	5	U
-Trimethylbenzene	95-63-6	5	U
ylbenzene	135-98-8	5	U
ichlorobenzene	541-73-1	5	U
propyltoluene	99-87-6	5	U
ichlorobenzene	106-46-7	5	U
ylbenzene	104-51-8	5	U
ichlorobenzene	95-50-1	5	U
ibromo-3-chloropropane	96-12-8	5	U
-Trichlorobenzene	120-82-1	5	U
chlorobutadiene	87-68-3	5	U
halene	91-20-3	10	U
-Trichlorobenzene	87-61-6	5	U

Gate Compound	%Rec	Recovery Limits (%)
Dibromofluoromethane	100 %	86 --- 118
Toluene-d8	103 %	88 --- 110
-Bromofluorobenzene	101 %	86 --- 116

## Qualifiers:

- "U" Indicates compound was searched for and not detected.
- "B" Indicates compound was found in the method blank.
- "J" Indicates compound was identified out of the method working limits and should be considered an estimated value.
- "D" Indicates compound was run at a dilution.
- "\*" Indicates surrogate recovery is not within method limits due to matrix effect.

Note: Method detection limits are approximately 1/5 of reporting limits.

## Page 3 of 3

Client Sample ID

Data Filename: VOAA3238.D

Units:  $\mu\text{g/L}$ 

**Qualifier:**

All tentatively identified compounds are estimated values.

**Analyst:**

Александр В. Яценко

# Petroleum Hydrocarbons Report

Page 1 of 1

3026-010; TH-25

ORIGINAL

PA Method: mod. 8015/mod. 8100  
Lab Sample ID: 99-3-30-2  
Matrix: Water  
Tag Number: 58999, 59000, 59001  
Date Sampled: 03/23/99

Analyst: DPD  
Volatiles Date Analyzed: 03/24/99  
Date Extracted: 03/24/99  
Extractables Date Analyzed: 03/24/99  
Units:  $\mu\text{g/L}$

Volatiles Dilution Factor: 1

Extractables Dilution Factor: 1

Analyte	CAS Number	Concentration	Reporting Limits	Qualifier
Total Volatile Hydrocarbons	NA		500	U
Total Extractable Hydrocarbons	NA		1000	U

Surrogate Compound	%Recovery
(SS) a,a,a-Trifluorotoluene	57 %
(SS) Fluorobenzene	102 %
(SS) o-Terphenyl	52 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank and has been corrected.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Dongming Qiu*



No. 8763

4888 Pearl East Circle, Suite 108  
Boulder Colorado 80301

PM: STAN SPENCER

99-3-30

[illegible]

# 3026-010; TH-26

Client Sample ID

Sample Tag No: 58390, 91, 92

Lab Sample ID: W-99-6-42-1

Matrix: Water

Date Sampled: 6/23/99

Data Filename: VOAA3432.D

VOAA3450.D

EPA Method: 8260

Date Analyzed: 6/23,24/99

Analyst: SBS

Units: µg/L

Dilution Factor: 1

2

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
1,1-Difluoromethane	75-71-8		20	DU
1,1-Dichloromethane	74-87-3		10	U
1,1,1-Trichloroethane	75-01-4		5	U
1,1,2-Trichloroethane	74-97-5		10	U
1,1,2,2-Tetrachloroethane	75-00-3		10	U
1,1,2,2-Tetrachloroethane	75-69-4		10	U
1,1-Dichloroethane	75-35-4		5	U
1,1,1-Trichloroethane	75-09-2		5	U
1,2-Dichloroethane	156-60-5		5	U
1,2-Dichloroethane	75-34-3		5	U
1,2-Dichloroethane	156-59-2		5	U
1,2-Dichloroethane	594-20-7		5	U
1,2-Dichloroethane	74-97-5		5	U
1,2-Dichloroethane	67-66-3		5	U
1,2-Dichloroethane	71-55-6		5	U
1,2-Dichloroethane	563-58-6		5	U
1,2-Dichloroethane	56-23-5		5	U
1,2-Dichloroethane	107-06-2		5	U
1,2-Dichloroethane	71-43-2		5	U
1,2-Dichloroethane	79-01-6		5	U
1,2-Dichloroethane	78-87-5		5	U
1,2-Dichloroethane	74-95-3		5	U
1,2-Dichloroethane	75-27-4		5	U
1,2-Dichloroethane	10061-02-6		5	U
1,2-Dichloroethane	108-88-3		5	U
1,2-Dichloroethane	10061-01-5		5	U
1,2-Dichloroethane	79-00-5		5	U
1,2-Dichloroethane	106-93-4		5	U
1,2-Dichloroethane	142-28-9		5	U
1,2-Dichloroethane	127-18-4	300	10	D
1,2-Dichloroethane	124-48-1		5	U
1,2-Dichloroethane	108-90-7		5	U
1,2-Dichloroethane	630-20-6		5	U
1,2-Dichloroethane	100-41-4		5	U
1,2-Dichloroethane			5	U
1,2-Dichloroethane	106-42-3		5	U



Walsh Environmental Scientists and Engineers, Inc.



3026-010; TH-26

Lab Sample ID: W-99-6-42-1

Client Sample ID:

ene	100-42-5		5	U
noform	75-25-2		5	U
rophenylbenzene	98-82-8		5	U
2,2-Tetrachloroethane	79-34-5		5	U
nobenzene	108-86-1		5	U
3-Trichloropropane	96-18-4		5	U
opphenylbenzene	103-65-1		5	U
lorotoluene	95-49-8		5	U
3-Trimethylbenzene	108-67-8		5	U
tylbenzene	98-06-6		5	U
4-Trimethylbenzene	95-63-6		5	U
tylbenzene	135-98-8		5	U
Dichlorobenzene	541-73-1		5	U
propyltoluene	99-87-6		5	U
Dichlorobenzene	106-46-7		5	U
tylbenzene	104-51-8		5	U
Dichlorobenzene	95-50-1		5	U
Dibromo-3-chloropropane	96-12-8		5	U
-Trichlorobenzene	120-82-1		5	U
chlorobutadiene	87-68-3		5	U
thalene	91-20-3		5	U
-Trichlorobenzene	87-61-6		5	U

rogate Compound	%Rec	Limits (%)
Dibromofluoromethane	85 %	70 --- 130
Toluene-d8	84 %	70 --- 130
p-Bromofluorobenzene	83 %	70 --- 130

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

\*\*\* Indicates surrogates low due to matrix effect.

"D" Indicates result obtained from a dilution.



Environmental Scientists and Engineers, Inc.

3026-010; TH-26

Client Sample ID

Sample Tag No: 58390, 91, 92

Lab Sample ID: W-99-6-42-1

Matrix: Water

Date Sampled: 06/23/99

Data Filename: VOAA3432.D

VOAA3450.D

EPA Method: 8260

Date Analyzed: 6/23,24/9

Analyst: SBS

Units:  $\mu\text{g/L}$ 

Dilution Factor: 1

2

[illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

**Analyst:**

Leifur K. Olsen



Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

ORIGINAL

3026-010; Brighton TH-26

Client Sample ID

Method: Mod. 8100  
Sample ID: 99-6-42-1  
Water  
Number: 58390/91/92  
Sampled: 6/23/99

Analyst: SBS  
Date Extracted: 6/23/99  
Extractables Date Analyzed: 7/2/99  
Units: mg/L

Extractables Dilution Factor: 1

	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		1	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	51 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"\*" Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: 

**Walsh**

Environmental Scientists and Engineers, Inc.

**3026-010; TH-27**

Client Sample ID

Sample Tag No: 58393. 94, 95

Lab Sample ID: W-99-6-42-2

Matrix: Water

Date Sampled: 6/23/99

Data Filename: VOAA3433.D

VOAA3449.D

EPA Method: 8260

Date Analyzed: 6/23,24/9

Analyst: SBS

Units: µg/L

Dilution Factor: 1

2

yte	CAS Number	Concentration	Quantitation Limits	Qualifier
orodifluoromethane	75-71-8		20	DU
omethane	74-87-3		10	U
Chloride	75-01-4		5	U
omethane	74-97-5		10	U
oethane	75-00-3		10	U
orofluoromethane	75-69-4		10	U
ichloroethene	75-35-4		5	U
/lene Chloride	75-09-2		5	U
1,2-Dichloroethene	156-60-5		5	U
ichloroethane	75-34-3		5	U
2-Dichloroethene	156-59-2		5	U
chloropropane	594-20-7		5	U
ochloromethane	74-97-5		5	U
oform	67-66-3		5	U
Trichloroethane	71-55-6		5	U
chloropropene	563-58-6		5	U
n Tetrachloride	56-23-5		5	U
chloroethane	107-06-2		5	U
ne	71-43-2		5	U
roethene	79-01-6		5	U
chloropropane	78-87-5		5	U
nomethane	74-95-3		5	U
odichloromethane	75-27-4		5	U
-Dichloropropene	10061-02-6		5	U
ie	108-88-3		5	U
,3-Dichloropropene	10061-01-5		5	U
Trichloroethane	79-00-5		5	U
romoethane	106-93-4		5	U
chloropropane	142-28-9		5	U
loroethene	127-18-4	340	10	D
ochloromethane	124-48-1		5	U
benzene	108-90-7		5	U
-Tetrachloroethane	630-20-6		5	U
enzene	100-41-4		5	U
Xylenes			5	U
ne	106-42-3		5	U

**Walsh**

ental Scientists and Engineers, Inc.



3026-010; TH-27

Lab Sample ID: W-99-6-42-2

Client Sample ID:

ene	100-42-5		5	U
noform	75-25-2		5	U
propylbenzene	98-82-8		5	U
2,2-Tetrachloroethane	79-34-5		5	U
nobenzene	108-86-1		5	U
3-Trichloropropane	96-18-4		5	U
propylbenzene	103-65-1		5	U
chlorotoluene	95-49-8		5	U
3-Trimethylbenzene	108-67-8		5	U
ylbenzene	98-06-6		5	U
-Trimethylbenzene	95-63-6		5	U
tylbenzene	135-98-8		5	U
Dichlorobenzene	541-73-1		5	U
propyltoluene	99-87-6		5	U
Dichlorobenzene	106-46-7		5	U
ylbenzene	104-51-8		5	U
Dichlorobenzene	95-50-1		5	U
ibromo-3-chloropropane	96-12-8		5	U
-Trichlorobenzene	120-82-1		5	U
chlorobutadiene	87-68-3		5	U
thalene	91-20-3		5	U
-Trichlorobenzene	87-61-6		5	U

rogate Compound	%Rec	Limits (%)
Dibromofluoromethane	83 %	70 --- 130
Toluene-d8	83 %	70 --- 130
p-Bromofluorobenzene	85 %	70 --- 130

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

"" Indicates surrogates low due to matrix effect.

"D" Indicates result obtained from a dilution.



mental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; Brighton TH-27**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-42-2  
ix: Water  
Number: 58393/94/95  
Sampled: 6/23/99

Analyst: SBS  
Date Extracted: 6/23/99  
Extractables Date Analyzed: 6/24/99  
Units: mg/L

Extractables Dilution Factor: 1.25

yte	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		1.25	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	91 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

*Stephen B. B. B.*

**Walsh**

Environmental Scientists and Engineers, Inc.

**3026-010; TH-28**

Client Sample ID

Sample Tag No: 58396. 97, 98

Lab Sample ID: W-99-6-42-3

Matrix: Water

Date Sampled: 6/23/99

Data Filename: VOAA3448.D

EPA Method: 8260

Date Analyzed: 06/24/99

Analyst: SBS

Units: µg/L

Dilution Factor: 1

Compound	CAS Number	Concentration	Quantitation Limits	Qualifier
Perdifluoromethane	75-71-8		10	U
Fluoromethane	74-87-3		10	U
Chloride	75-01-4		5	U
Fluoromethane	74-97-5		10	U
Methane	75-00-3		10	U
Perfluoromethane	75-69-4		10	U
Chloroethene	75-35-4		5	U
Ethene Chloride	75-09-2		5	U
1,2-Dichloroethene	156-60-5		5	U
Chloroethane	75-34-3		5	U
1,1-Dichloroethene	156-59-2		5	U
Chloropropane	594-20-7		5	U
Chloromethane	74-97-5		5	U
Form	67-66-3		5	U
Trichloroethane	71-55-6		5	U
Chloropropene	563-58-6		5	U
1,1,1-Tetrachloride	56-23-5		5	U
Chloroethane	107-06-2		5	U
Ethene	71-43-2		5	U
Chloroethene	79-01-6		5	U
Chloropropane	78-87-5		5	U
Fluoromethane	74-95-3		5	U
Dichloromethane	75-27-4		5	U
Dichloropropene	10061-02-6		5	U
Ethene	108-88-3		5	U
1,3-Dichloropropene	10061-01-5		5	U
Trichloroethane	79-00-5		5	U
Fluoroethane	106-93-4		5	U
Chloropropane	142-28-9		5	U
Chloroethene	127-18-4	160	5	
Dichloromethane	124-48-1		5	U
Benzene	108-90-7		5	U
Tetrachloroethane	630-20-6		5	U
Benzene	100-41-4		5	U
Cylenes			5	U
Ethene	106-42-3		5	U

**Walsh**

Environmental Scientists and Engineers, Inc.



3026-010; TH-28

Client Sample ID

Sample Tag No: 58396, 97, 98

Lab Sample ID: W-99-6-42-3

**Matrix:** Water

Date Sampled: 06/23/99

Data Filename: VOAA3448.D

EPA Method: 8260

Date Analyzed: 06/24/99

Analyst: SBS

Units:  $\mu\text{g/L}$ 

Dilution Factor: 1

[illegible]

**Qualifier:**

"T" Indicates compound was tentatively identified by its mass spectrum.

All tentatively identified compounds are estimated values.

**Analyst:**

Die gute Nacht



**mental Scientists and Engineers, Inc.**

3026-010; TH-28

Lab Sample ID: W-99-6-42-3

Client Sample ID:

Styrene	100-42-5		5	U
Bromoform	75-25-2		5	U
Isopropylbenzene	98-82-8		5	U
1,1,2,2-Tetrachloroethane	79-34-5		5	U
Bromobenzene	108-86-1		5	U
1,2,3-Trichloropropane	96-18-4		5	U
n-Propylbenzene	103-65-1		5	U
2-Chlorotoluene	95-49-8		5	U
1,3,5-Trimethylbenzene	108-67-8		5	U
t-Butylbenzene	98-06-6		5	U
1,2,4-Trimethylbenzene	95-63-6		5	U
s-Butylbenzene	135-98-8		5	U
1,3-Dichlorobenzene	541-73-1		5	U
p-Isopropyltoluene	99-87-6		5	U
1,4-Dichlorobenzene	106-46-7		5	U
n-Butylbenzene	104-51-8		5	U
1,2-Dichlorobenzene	95-50-1		5	U
1,2-Dibromo-3-chloropropane	96-12-8		5	U
1,2,4-Trichlorobenzene	120-82-1		5	U
Hexachlorobutadiene	87-68-3		5	U
Naphthalene	91-20-3		5	U
1,2,3-Trichlorobenzene	87-61-6		5	U

Surrogate Compound	%Rec	Limits (%)
(SS) Dibromofluoromethane	74 %	70 -- 130
(SS) Toluene-d8	77 %	70 -- 130
(SS) p-Bromofluorobenzene	74 %	70 -- 130

Qualifiers: "U" Indicates compound was searched for and not detected.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

\*\*\* Indicates surrogates low due to matrix effect.



Environmental Scientists and Engineers, Inc.

# Petroleum Hydrocarbons Report

Page 1 of 1

**3026-010; Brighton TH-28**

Client Sample ID

**ORIGINAL**

Method: Mod. 8100  
Sample ID: 99-6-42-3  
Matrix: Water  
Number: 58396/97/98  
Sampled: 6/23/99

Analyst: SBS  
Date Extracted: 6/23/99  
Extractables Date Analyzed: 6/24/99  
Units: mg/L

Extractables Dilution Factor: 1

Compound	CAS Number	Concentration	Detection Limits	Qualifier
Extractable Hydrocarbons	NA		1	U

Surrogate Compound	%Recovery
(SS) o-Terphenyl	51 %

## Qualifiers:

"U" Indicates compound was searched for and not detected at or above the method detection limit.

"B" Indicates compound was found in the method blank.

"J" Indicates compound was identified out of the method working limits and should be considered an estimated value.

" \* " Indicates surrogate is outside of recovery limits due to matrix effect.

Analyst: \_\_\_\_\_

**Valsh**

Scientists and Engineers, Inc.

480546

# Superfund DOCUMENT CODING SHEET

SITE NAME (Required): VASQUEZ BOULEVARD AND I-70

CERCLIS NUMBER: C00002259588

SSID: (Add the two digit Site Spill ID #) 08 - 9R

OPERABLE UNIT (Add the applicable operable unit(s): 01 and 02

PHASE ACTIVITY (Required): (Check one or more boxes)

- |   |  |   |  |
|---|--|---|--|
| <input type="checkbox"/> Brownfields                                | <input type="checkbox"/> Natural Resource Damages  | <input type="checkbox"/> Post Work Cost Recovery Action   | <input checked="" type="checkbox"/> Remedial Studies/ Remy Selection |
| <input type="checkbox"/> Emergency Response/ Time Critical Removals | <input type="checkbox"/> Non-Time Critical Removal | <input type="checkbox"/> Remedial Design/ Remedial Action | <input type="checkbox"/> Removal Site Evaluation                     |
| <input type="checkbox"/> HRS  | <input type="checkbox"/> Post Construction         |   | <input type="checkbox"/> Site Assessment                             |

DOCUMENT TYPE (Required): (Check one or more boxes)

- |   |  |  |  |
|---|--|--|--|
| <input type="checkbox"/> Access Related                     | <input type="checkbox"/> Correspondence Internal       | <input type="checkbox"/> Financial & Contract Documents            | <input type="checkbox"/> Orders, Settlements                 |
| <input type="checkbox"/> CERCLIS Documentation              | <input checked="" type="checkbox"/> Correspondence PRP | <input type="checkbox"/> HRS Package                               | <input type="checkbox"/> Plans (Workplans, QAPPs, CRP, HASP) |
| <input type="checkbox"/> Comfort Letters                    | <input type="checkbox"/> Data & Related Documents      | <input type="checkbox"/> IAGs, MOAs, SSCs & CAs                    | <input type="checkbox"/> Reports                             |
| <input type="checkbox"/> Community Relations                | <input type="checkbox"/> Decision Documents            | <input type="checkbox"/> Map (Photo, Video, Tape, etc)             | <input type="checkbox"/> Risk Assessment - Ecological        |
| <input type="checkbox"/> Construction Related               | <input type="checkbox"/> Design Related                | <input checked="" type="checkbox"/> Notice Letter & 104(e) Related | <input type="checkbox"/> Risk Assessment - Human Health      |
| <input checked="" type="checkbox"/> Correspondence External | <input type="checkbox"/> Enforcement                   |  |  |

REFERENCE: (List PRPs Document Pertains to Below)

- ☒ DENVER
- ☐
- ☐
- ☐
- ☐
- ☐
- ☐

ACCESS CODE (Required): (Check Applicable Boxes)

- |  |   |
|--|---|
| <input type="checkbox"/> Attorney Work Product | <input type="checkbox"/> Deliberative Process           |
| <input type="checkbox"/> Attorney-Client Comm. | <input type="checkbox"/> OGC OK                         |
| <input type="checkbox"/> CBI Claimed           | <input type="checkbox"/> Enforcement Confidential       |
| <input type="checkbox"/> CBI Determined        | <input type="checkbox"/> Privacy                        |
|  | <input checked="" type="checkbox"/> Releasable (Public) |

RELATED DOCUMENTS (ATTACHMENTS):

Total Number of Attachment(s): \_\_\_\_\_

Title of Attachment(s): \_\_\_\_\_

Number of Attachment(s): \_\_\_\_\_

OF \_\_\_\_\_

OF \_\_\_\_\_

OF \_\_\_\_\_

OF \_\_\_\_\_

COLLECTIONS: (Write the title of your collection on the line provided. Then check the box to add the document to the selected collection.)

- | Administrative Record Collections                          | Special Collections                     | FOIA Collections         | Discovery Collections    | Cost Recovery Collections |
|--|---|--------------------------|--------------------------|---------------------------|
| <input checked="" type="checkbox"/> <u>104(e) response</u> | <input type="checkbox"/> Work Performed | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>  |
| <input type="checkbox"/> <u>from Denver</u>                | <input type="checkbox"/>                | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>  |
| <input type="checkbox"/>                                   | <input type="checkbox"/>                | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>  |
| <input type="checkbox"/>                                   | <input type="checkbox"/>                | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/>  |

NOTES TO RECORDS CENTER:

Please list attachments separately.

SIGNATURE BOX (Required)

Program Lead: Mark Herman

Attorney: Mark Herman

DATE: 4/27/00

DATE: 11/15/00